

GENERAL ATOMIC DIVISION
CENTRAL DYNAMICS CORPORATION

John Jay Hopkins Laboratory for Pure and Applied Science
P.O. Box 600, San Diego 12, California

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A CONTINUOUS TWO-DIMENSIONAL EULERIAN HYDRODYNAMIC CODE

Work done by:

Project Staff

Approved by:

W. J. Johnson

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FOREWORD

The GIL computer code described herein is as it existed on October 26, 1964. The code has been in continuous development for three years and in its presented form has been applied successfully by General Atomic to the kind of problems discussed in this report. However, the development and improvement in both the physics and mathematics of the code are being continued, so that duplication of results (or even close agreement) between problems run with the code as published and the code as it existed either before or after this time is not necessarily to be expected.

General Atomic has exercised due care in preparation, but does not warrant the merchantability, accuracy, and completeness of the code or of its description contained herein. The complexity of this kind of program precludes any guarantee to that effect. Therefore, any user must make his own determination of the suitability of the code for any specific use and of the validity of the information produced by use of the code.

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1. INTRODUCTION AND ACKNOWLEDGEMENTS

The OIL code is very closely related to familiar particle-in-cell codes^(1,2,3) and has been developed by modification of the General Atomic particle-in-cell code named SHELL. The basic difference lies in the method of the mass transport, the OIL scheme being a continuous mass transport, while the SHELL scheme is discrete. The initial work on a continuous version of SHELL was undertaken several years ago by B. E. Freeman and the author. Since early 1963 the development of the continuous Eulerian has been continued by J. M. Walsh and the author.

The author is deeply appreciative of the work and consultation given by B. E. Freeman and J. M. Walsh. The assistance of D. R. Yates and R. H. Fisher in the automatic plotting routines used for the OIL code is also much appreciated.

Since its original use for the hypervelocity impact calculations⁽⁴⁾, the OIL code has been successfully adapted to several other high energy fluid dynamic applications.

Detailed descriptions of various problems, especially results for hypervelocity impact, are given in Ref. 4.

2. CLAM

2.1. General Description

CLAM is the generator code for the OIL code, and is used to generate initial values of the variables for each cell in the grid. (An exception arises for certain simple initial conditions, described in Section 3.2, where it is possible to bypass CLAM and use instead a more economical routine called SETUP.)

The function of CLAM is illustrated by a simple example (Fig. 1). We wish to generate the following grid: A right circular cylinder of density $1. \text{ g/cm}^3$, radius 24 cm, and height 12 cm is located at position $z = 20 \text{ cm}$ along the axis. A projectile (right cylinder) of density $1. \text{ g/cm}^3$, radius 6 cm, and height 12 cm, is located at positions 8-20 along the axis. The projectile has an initial axial velocity of $1 \times 10^6 \text{ cm/sec}$. The cells are 1 cm on a side.

2

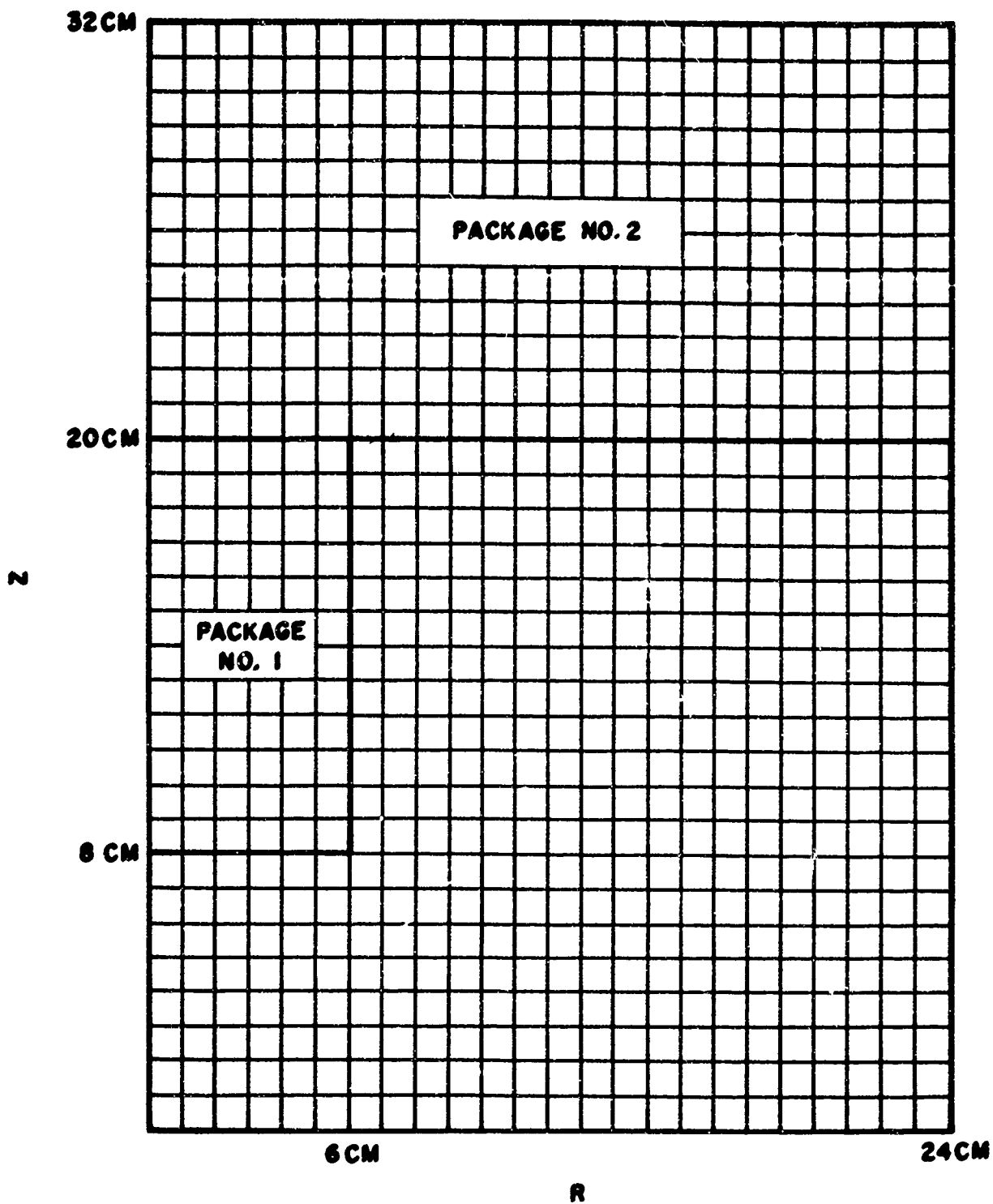


Fig. 1

3.

(CLAM INPUT)

		GAS	GAMMA = 1.5	RHO TARGET =	RHO PELLET = 1.		
1.		24.	32.	0.	2.	0.	07
0024		1.					
1132		1.					
2.		3.	80.				
11	01	0.	0.	1.			
4	1	0.	6.	8.	20.		
51		1-0	0.	0.	0.	0.	0.
52		0.	0.	0.	0.	0.	0.
53		.01	0.	0.	0.	0.	0.
11	01	0.	0.	1.			
4	1	0.	24.	20.	32.		
53		0.	0.	0.	0.	0.	0.
2							

GAS GAMMA = 1.5 RHO TARGET= RHO PELLET = 1.

PROB. NO. 1.000

I=24

J=30

X(I) I=0,24

0.0	1.0	2.0	3.0	4.0
5.0	6.0	7.0	8.0	9.0
10.0	11.0	12.0	13.0	14.0
15.0	16.0	17.0	18.0	19.0
20.0	21.0	22.0	23.0	24.0

Y(J) J=0,32

0.0	1.0	2.0	3.0	4.0
5.0	6.0	7.0	8.0	9.0
10.0	11.0	12.0	13.0	14.0
15.0	16.0	17.0	18.0	19.0
20.0	21.0	22.0	23.0	24.0
25.0	26.0	27.0	28.0	29.0
30.0	31.	32.0		

DX(I) I=1,24

1.0	1.0	1.0	1.0	1.0
1.0	1.0	1.0	1.0	1.0
1.0	1.0	1.0	1.0	1.0
1.0	1.0	1.0	1.0	1.0
1.0	1.0	1.0	1.0	1.0

4.

DY(J) J=1,32.

1.0	1.0	1.0	1.0	1.0
1.0	1.0	1.0	1.0	1.0
1.0	1.0	1.0	1.0	1.0
1.0	1.0	1.0	1.0	1.0
1.0	1.0	1.0	1.0	1.0
1.0	1.0	1.0	1.0	1.0
1.0	1.0	1.0	1.0	1.0
1.0	1.0	1.0	1.0	1.0

AREA(I) I=1,48

3.141593	9.424778	15.707963	21.991149	28.274334
34.557519	40.840705	47.123890	53.407076	59.690261
65.973447	72.256632	78.539817	84.823003	91.106188
97.389374	103.672559	109.955745	116.238930	122.522116
128.805300	135.069486	141.371672	147.654856	

PACKAGE NO. 1 1 PARTICLE /CELL

	A1	A2	A3	A4	A5	A6
DENSITY	1.	0.	0.	0.	0.	0.
ENERGY	0.	0.	0.	0.	0.	0.
VELOCITY	0.01	0.	0.	0.	0.	0.

RECTANGLE---GEN--- 0. 6. 8. 20.
I(1)= 1 J(1)= 8 I(N)= 7 J(N)= 21

72 (X) PARTICLES PE=6.785839-02 PM=1.357168+03

PACKAGE NO. 2 1 PARTICLE /CELL

	A1	A2	A3	A4	A5	A6
VELOCITY	0.	0.	0.	0.	0.	0.

RECTANGLE--- GEN--- 0. 24. 20. 32.
I(1)= 1 J(1)=20 I(N)=24 J(N)=32

288 (X) PARTICLES PE=0. PM =2.171467+04

THE = 6.785838399-2 E = 6.785839237-2

M. = 0. MX = 2.30718+04 M.+ MX = 2.30718+04

PARTICLES --- 0 DOT 360 X 360 TOTAL

TAPE DUMP AT TIME 0. CYCLE 0.

The coordinates $x(i)$ and $y(i)$ of each cell are calculated from the DX 's and DY 's that are loaded. In addition, the area term $(TAU_{(i)}) = (x_{(i)}^2 - x_{(i-1)}^2)\pi$ are calculated.

The next step is to process the packages (see Section 2.2); there are two in this example. Package Number (1) is the projectile and Package Number (2) is the target.

We specify the coordinates of each package. In addition, we specify the number of particles per cell in each package. We assign a density, the two velocity components, and specific internal energy for the package in question. These may be constants or functions of both z and r .

After specifying the coordinates of the package, particles (N , where N may range from 1 to 20) are placed in each cell. Each of the particles is then assigned a density, the two velocity components and specific internal energy. The cell is divided into N equal parts, and the N particles placed at the center of these areas. The mass of each particle is then the density (prescribed along with the package input) times the volume of the small subdivision cell of the cell (K). The mass of cell (K) then is the sum of the N particle masses, and both momenta components are calculated as the sum of the individual momenta components of each particle. The internal energy of each cell (K) is the specific internal energy times the particle mass summed over all N particles.

In the output routine these cell quantities are then converted to the two velocity components and the specific internal energy.

Particles are created, whether these data are for an OIL or a SHELL run, for the purpose of describing the many possible geometries and possible energy, velocity, and density distributions. If this is an OIL run, the particles are not written on a tape as they would be if this were a SHELL run. One continues to process each cell (K) within the given package and the N particles in cell (K).

In this example, it is sufficient to specify only one particle per cell, since the package boundaries coincide with the cell boundaries and the density, velocities, and internal energy are constant.

After processing all packages, one then writes this information about the grid and the cells on a tape that will be the starting conditions for the OIL or SHELL code (see Section 2.3.)

The output from CLAM, excluding the long print, is shown on the listings on pages 3 and 4.

2.2. Input Description for CLAM

CLAM and OIL are written in cylindrical coordinates with axial symmetry. In the following discussion and description, X refers to the coordinate R and Y to the coordinate Z. An asterisk before the word signifies that the data are in floating point; otherwise they are fixed point.

<u>Card No.</u>	<u>Column No.</u>	<u>Description</u>
1		Header card, any BCD information in col. 2-72.
2	* 1-10	Contain the problem number. If less than zero, this will be a particle run using SHELL; if greater than zero, this will be a continuous run using OIL.
	* 11-20	IMAX, the number of cells in the X-direction (maximum of 50.)
	* 21-30	JMAX, the number of cells in the Y-direction (maximum of 100)
	* 31-40	Blank
	* 41-50	2
	* 51-60	S8 (not used as input in CLAM)
	* 61-70	S9 (not used in CLAM)
	* 71-72	N7 (N7 = binary tape number)
3		(2 cards is the minimum)
	1	A(1) indicates that this is the last DX or DY card to be read in. A(0) indicates that there will be more DX or DY data cards.
	2	A(0) indicates that we are loading DX data A(1) indicates DY data.
	3-4	Indicate the number of zones that will have the same DX or DY values that appear in columns 11-20.
	5-6	Indicate the number of zones that will have the same DX or DY values that appear in columns 21-30.

<u>Card No.</u>	<u>Column No.</u>	<u>Description</u>
3	7-8	Indicate the number of zones that will have the same DX or DY values that appear in columns 31-40.
	9-10	Indicate the number of zones that will have the same DX or DY values that appear in columns 41-50.
	* 11-20	The value of DX or DY
	* 21-30	The value of DX or DY
	* 31-40	The value of DX or DY
4	* 41-50	The value of DX or DY
	* 1-10	N1 = the tape number of one of the scratch tapes to be used in CLAM and OIL
	* 11-20	N2 = the other tape number for the scratch tape. OIL requires two scratch tapes if using the particle transport.
	* 21-30	N4 = maximum number of particles plus one per tape record, that CLAM will generate (maximum value = 130)
	* 31-40	Switch (not used)

Now we begin loading the necessary data to generate a package. The maximum number of packages that may be generated is 72; to increase the maximum requires changing the dimension statements.

1	1	Load a 1 here
	2	A(1) implies that x material will be generated in this package. A(0) implies that dot material will be generated.
	5-7	(N ²), the number of particles per cell to be generated, where $1 \leq N \leq 20$.
	* 11-20	YC = Y coordinate for the origin of the radius vector used in the density, energy, and velocity fits.
	* 21-30	XC = X coordinate for the origin of the radius vector used in the density, energy, and velocity fits.
	* 31-40	BMP(3) is loaded into S8 (S8 contains the number (1-6) of the fit for the density, internal energy, and velocity that will be assigned to each particle as it is generated in this package.
	* 41-70	Blank.

Following the first card of each package are five other types of cards:

- (1) Generate (geometry of package)
- (2) Delete card (if needed); there may be more than one
- (3) A density card (only one per package)
- (4) An energy card (only one per package)
- (5) A velocity card (only one per package)

For card (1) above, CLAM has the following geometric options for generating or deleting:

1. A rectangle - a 4 in column 1
 - a (1) in column 7 means generate this rectangle
 - a (0) in column 7 means delete this rectangle
 - * 11-20 X1 = the left R coordinate of the rectangle
 - * 21-30 X2 = the right R coordinate of the rectangle
 - * 31-40 Y1 = the lower Z coordinate of the rectangle
 - * 41-50 Y2 = the upper Z coordinate of the rectangle
2. A triangle - a 6 in column 1
 - a (1) in column 7 means generate this triangle
 - a (0) in column 7 means delete this triangle.
 - * 11-20 X1)
 - * 21-30 Y1)
 - * 31-40 X2) --Note: Vertices (1-3) can be in any order
 - * 41-50 Y2)
 - * 51-60 X3)
 - * 61-70 Y3)
3. An ellipse or circle - a 41 in column (1-2)
 - a (1) in column 7 means generate this ellipse or circle
 - a (0) in column 7 means delete this ellipse or circle

- * 11-20 The semi-axis in the X-direction if an ellipse or the radii if for a circle
- * 21-30 The semi-axis in the Y-direction if an ellipse or zero if for a circle
- * 31-40 The X-coordinate of the center of ellipse or circle
- * 41-50 The Y-coordinate of the center of ellipse or circle.

4. A perturbed ellipse - a 61 in columns (1-2)

a (1) in column 7 means generate this perturbed ellipse.

a (0) in column 7 means delete this perturbed ellipse.

- * 11-20 Semi-axis in the X-direction
- * 21-30 Semi-axis in the Y-direction
- * 31-40 0.
- * 41-50 Y-coordinate of center of perturbed ellipse
- * 51-60 X-coordinate of point of perturbation
- * 61-70 Y-coordinate of point of perturbation.

Following the geometry cards are the following:

Density card - a 51 in columns (1-2)

Energy card - a 52 in columns (1-2)

Velocity card - a 53 in columns (1-2)

Note: If in this package, the ρ or I or velocity will remain the same as the previous package, then a 51, 52 or 53 card is not required.

- * 11-20)
- * 21-30)
- * 31-40) - Contains the values to be used in the analytical
- * 41-50) expressions for the density, energy, and velocities.
- * 51-60)
- * 61-70)

This data is then loaded into the following arrays:

TABR(1-6) for density

TABI(1-6) for internal energy

TABUV(1-6) for velocity

Finally, a 2 in column 1 signifies the completion of loading all input cards for the CLAM code.

There are six subroutines (FIT1 - FIT6) used for computing ρ , I, U, and V. The standard input to these subroutines is as follows:

TY = Y coordinate of particle N

TX = X coordinate of particle N

The modified coordinates TTY and TTX are computed as follows:

TTY = Y coordinate = TY - YC (relative to YC)

TTX = X coordinate = TX - XC (relative to XC)

The standard output from the subroutines is as follows:

WSR = ρ (density) of particle N

WSI = I (specific internal energy) of particle N

WSU = U (radial velocity component) of particle N

WSV = V (axial velocity component) of particle N.

Below are the equations, or analytical fits, for the six subroutines. Any or all may be changed. Each equation is followed by the FORTRAN mnemonic.

1. FIT 1

$$R = (X^2 + Y^2)^{\frac{1}{2}}$$

$$WS = (TTX^2 + TTY^2)^{\frac{1}{2}}$$

$$\rho = A + B (Y - C)$$

$$WSR = TABR(1) + TABR(2) * (TTY - TABR(3))$$

$$I = A + B (Y - C)$$

$$WSI = TABI(1) + TABI(2) * (TTY - TABI(3))$$

$$U = 0$$

$$WSU = 0$$

$$V = A + B (Y - C)$$

$$WSV = TABUV(1) + TABUV(2) * (TTY - TABUV(3))$$

2. FIT 2

$$R = (X^2 + Y^2)^{\frac{1}{2}}$$

$$WS = (TTX^2 + TTY^2)^{\frac{1}{2}}$$

$$\rho = \left(\frac{X-A}{B}\right)^2 + \left(\frac{Y-C}{D}\right)^2$$

$$WSR = \left(\frac{TTX - TABR(1)}{TABR(2)} \right)^2 + \left(\frac{TTY - TABR(3)}{TABR(4)} \right)^2$$

$$I = A + BX + CX^2 + DY + EY^2$$

$$WS1 = TAB1(1) + TAB1(2) * TTX + TAB1(3) * TTX^2 \\ + TAB1(4) * TTY + TAB1(5) * TTY^2$$

$$U = C + D * Y$$

$$WSU = TABUV(3) + TABUV(4) * TTY$$

$$V = A + B * Y$$

$$WSB = TABUV(1) + TABUV(2) * TTY$$

3. FIT 3

$$R = (X^2 + Y^2)^{\frac{1}{2}}$$

$$WS = (TTX^2 + TTY^2)^{\frac{1}{2}}$$

$$\rho = A + BR + CR^2$$

$$WSR = TABR(1) + TABR(2) * WS + TABR(3) * WS^2$$

$$I = A + BR + CR^2$$

$$WS1 = TAB1(1) + TAB1(2) * WS + TAB1(3) * WS^2$$

$$U = \frac{X}{R} \left(\frac{A + BR + CR^2}{D + ER + FR^2} \right)$$

$$WSU = \frac{TTX}{WS} \left(\frac{TABUV(1) + TABUV(2) * WS + TABUV(3) * WS^2}{TABUV(4) + TABUV(5) * WS + TABUV(6) * WS^2} \right)$$

$$V = \frac{Y}{R} \left(\frac{A + BR + CR^2}{D + ER + FR^2} \right)$$

$$WSV = \frac{TTY}{WS} \left(\frac{TABUV(1) + TABUV(2) * WS + TABUV(3) * WS^2}{TABUV(4) + TABUV(5) * WS + TABUV(6) * WS^2} \right)$$

Fits 4, 5, and 6 are dummy routines. Although particles are not used in the SHELL code if it is a continuous run (problem number greater than zero), the use of particles in CLAM provides the method for assigning mass, energy, to each cell.

2.3. OUTPUT FROM CLAM

The output from the CLAM code is the entire Z block (defined below), all the cell quantities (the two velocity components, the mass and internal energy), and the cell dimensions and areas. In the case where it is a particle run, the particles (their two coordinates and mass) and the i and j of the cell where the particle is located are also put onto the binary tape.

The normal system of units are the centimeter-gram-shake, where the units of energy are jerks/g and the pressure in units of jerks/cm³ (1 jerk = 10^{16} ergs and 1 shake = 10^{-8} sec).

The Z block or array contains the number of cells, the number of zones in both directions, and other necessary information to start the OIL or SHELL calculation. Below is a complete list of generated data from CLAM that is written on the binary output tape.

<u>Z</u>	<u>Equiv.</u>	<u>Units</u>	<u>Description</u>
1	PROB	--	Equals problem number, input to CLAM
2	Cycle	--	Equals cycle number = 0
3	DT	shake	Set to 0 by CLAM
4	Prints	--	↓
5	Print-L	--	
6	DUMPT 7	--	
7	C Stop	--	
8	PIDY	--	Equals $\pi = 3.1415927$
9	TMZ	grams	Total mass (x) generated by CLAM
10	GAM	--	Set to 0 by CLAM
11	GAMD	--	Set to 0 by CLAM
12	GAMX	--	Set to 0 by CLAM
13	ETH	jerk	Total energy in system
14	FFA	--	Set to 0 by CLAM
15	FFB	--	Set to 0 by CLAM
16	TMDZ	grams	Total mass (·) generated by CLAM; this version of CLAM does not generate (·)
17	TMXZ	grams	Total mass (x) generated by CLAM
18	XMAX	cm	= X(IMAX)

<u>Z</u>	<u>Equiv.</u>	<u>Units</u>	<u>Description</u>
19	TXMAX	cm	= 2 XMAX
20	TYMAX	cm	= 2. Y MAX (note Y MAX is not in Z block)
21	AMDM	grams	= minimum mass/2. of the dot particles
22	AMXM	grams	= minimum mass/2. of the X particles
23	DNN	--	Set to 0. by CLAM
24	DMIN	--	Set to 0. by CLAM
25	FEF	--	Set to 0. by CLAM
26	DTNA	--	Set to 0. by CLAM
27	CVIS	--	Set to 0. by CLAM
28	NPR	--	Set equal to 6 in CLAM
29	NPR1	--	CLAM sets NPR1 = N4 (check definition of N4(Z(54)))
30	NC	--	Fixed value of cycle number, set to 0. by CLAM
31	NPC	--	Used as indices in CLAM
32	NRC	--	Used as indices in CLAM
33	IMAX	--	Input to CLAM = maximum number of zones in X direction for this run
34	IMAX	--	Equal IMAX + 1
35	JMAX	--	Input to CLAM = maximum number of zones in Y direction for this run
36	JMAXA	--	= JMAX + 1
37	KMAX	--	= (IMAX)(JMAX) + 1
38	KMAXA	--	= KMAX + 1
39	NMAX	--	= total number of particles + 1 that CLAM has generated
40	ND	--	= total number of dot particles + 1 that CLAM has generated
41	KDT	--	Set to 0. by CLAM.
42	IXMAX	--	= IMAXA + 1
43	NOD	--	Used as index
44	NOPR	--	Set equal to N3 (Note definition of N3 (Z(53)))
45	NIMAX	--	Set to 0. by CLAM
46	NJMAX	--	Set to 0. by CLAM

<u>Z</u>	<u>Equiv.</u>	<u>Units</u>	<u>Description</u>
47	11		Set to 0. by CLAM
48	12		Set to 0. by CLAM
49	13		Set to 0. by CLAM
50	14	--	Set to 0. by CLAM
51	N1	--	= scratch tape number
52	N2	--	= scratch tape number
53	N3	--	= number of particle records of length $L_0 - 1$ that CLAM has generated
54	N4	--	= number of particles + 1. to be stored on each particle tape record
55	N5	-	Set to 0. by CLAM
56	N6	--	= number of particles on the last particle tape record
57	N7	--	= binary tape designation number
58	N8		Set to 0. by CLAM
59	N9		
60	N10		
61	N11		
62	NRM		
63	TRAD		
64	XNRG		
65	SN		
66	DXN		
67	RADER		
68	RADET		
69	RADEB		
70	DTRAD		
71	REZFCT		
72	RSTOP		
73	SHELL		<div style="text-align: center;">↓</div> A counter that may be used to distinguish between codes
74	BBOUND		Set to 0. by CLAM
75	TOZONE		Set to 0. by CLAM

<u>Z</u>	<u>Equiv.</u>	<u>Units</u>	<u>Description</u>
76	ECK		Set to 0. by CLAM
77	SBOUND		
78	X1		
79	X2		
80	Y1		
81	Y2		
82	CABLN		
83	VISC		
84	T		
85	GMAX		
86	WSGD		
87	WSGX		
88	GMADR		
89	GMAXR		
90	S1		
91	S2		
92	S3		
93	S4		
94	S5		
95	S6		
96	S7		
97	S8	--	Used for storage of FIT number for each package in CLAM
98	S9		Set to 0. by CLAM
99	S10		Set to 0. by CLAM

Z(100) through Z(150) is also set to 0. by CLAM.

3. OIL

3.1. The Eulerian equations we wish to solve are the following:

$$(A) \quad \frac{\partial \rho}{\partial t} + \nabla \cdot (\rho \vec{u}) = 0$$

$$(B) \quad \frac{\partial \rho \vec{u}}{\partial t} + \nabla \cdot (\rho \vec{u} \vec{u}) = -\nabla P$$

$$(C) \quad \frac{\partial \rho E}{\partial t} + \nabla \cdot (\rho E \vec{u}) = -\nabla \cdot (P \vec{u})$$

Equation (A) is the conservation of mass equation, (B) is the conservation of momentum, and (C) is the conservation of energy equation.

The second terms on the left side of Eqs. (B) and (C) are temporarily dropped. Their contributions are later approximated when we move particles or continuous mass across cell boundaries.

Rewriting equations (A), (B), and (C) in cylindrical coordinates with axis of symmetry results in Eqs. (1), (2), (3), and (4).

$$\frac{\partial \rho}{\partial t} = -\frac{\partial r \rho u}{r \partial r} - \frac{\partial \rho v}{\partial z} \quad (1)$$

$$\rho \frac{\partial u}{\partial t} = -\frac{\partial P}{\partial r} \quad (2)$$

$$\rho \frac{\partial v}{\partial t} = -\frac{\partial P}{\partial z} \quad (3)$$

$$\rho \frac{\partial E}{\partial t} = -\frac{\partial r P u}{r \partial r} - \frac{\partial P v}{\partial z} \quad (4)$$

$$P = f(\rho, I) \quad \text{Equation of state} \quad (5)$$

ρ = density of cell (K) in g/cm³

r = r coordinate in cm.

z = z coordinate in cm.

u = radial component of velocity in cm/shake

v = axial component of velocity in cm/shake

P = material pressure in jerks/cm³

E = total specific energy in jerks/g

I = specific internal energy in jerks/g

(1 jerk = 10¹⁶ ergs)

The five variables listed are all located at the center of the cell (Fig.2). For complete description of all quantities used, see Section 3.4 on List of Common for OIL."

Rewriting Eq. (4):

$$\rho \frac{\partial}{\partial t} [I + \frac{1}{2} (u^2 + v^2)] = - \frac{\partial r P u}{r \partial r} - \frac{\partial P v}{\partial z}$$

or

$$\rho \frac{\partial I}{\partial t} + \rho u \frac{\partial u}{\partial t} + \rho v \frac{\partial v}{\partial t} = - \frac{P}{r} \frac{\partial u r}{\partial r} - u \frac{r \partial P}{r \partial r} - v \frac{\partial P}{\partial z} - P \frac{\partial v}{\partial z}$$

but

$$\rho \frac{\partial u}{\partial t} = - \frac{\partial P}{\partial r} \quad \text{and} \quad \rho \frac{\partial v}{\partial t} = - \frac{\partial P}{\partial z}$$

thus

$$\rho \frac{\partial I}{\partial t} = - P \left(\frac{\partial v}{\partial z} + \frac{1}{r} \frac{\partial u r}{\partial r} \right)$$

Rewriting the momentum equations as

$$\rho \frac{\partial (u^2)}{\partial t} = - 2u \frac{\partial P}{\partial r} \quad \text{and} \quad \rho \frac{\partial (v^2)}{\partial t} = - 2v \frac{\partial P}{\partial z}$$

the radial momentum equation becomes in difference form

$$\rho \frac{\partial (u^2)}{\partial t} = 2u_{1-\frac{1}{2}, j-\frac{1}{2}} r_{1-\frac{1}{2}} \frac{[P_{1-3/2, j-\frac{1}{2}} - P_{1+\frac{1}{2}, j-\frac{1}{2}}]}{\Delta(r_1^2)}$$

and the axial momentum equation becomes

$$\rho \frac{\partial (v^2)}{\partial t} = v_{1-\frac{1}{2}, j-\frac{1}{2}} \frac{[P_{j-3/2, 1-\frac{1}{2}} - P_{j+\frac{1}{2}, 1-\frac{1}{2}}]}{\Delta z_j}$$

and the change in specific internal energy becomes

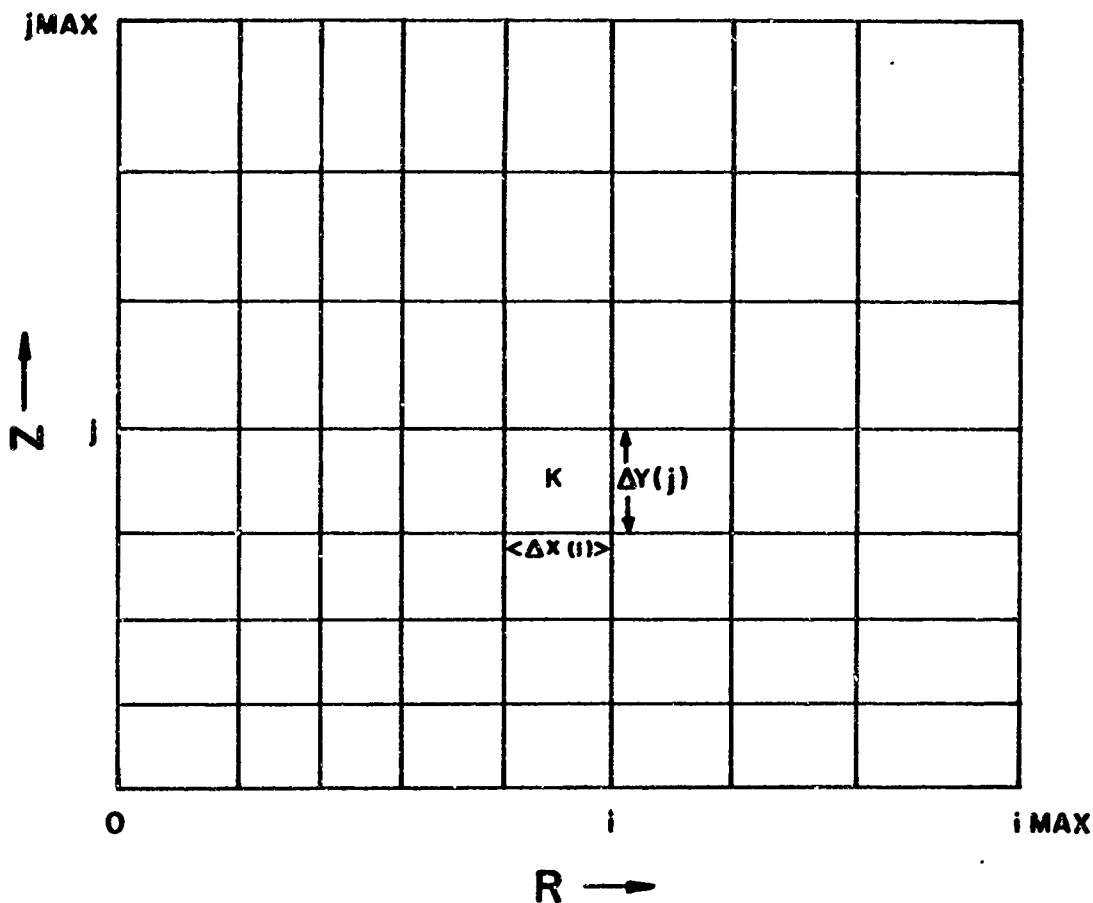
$$\rho \frac{\partial I}{\partial t} = - P \left[\frac{v_{1-\frac{1}{2}, j+\frac{1}{2}}^{n+\frac{1}{2}} - v_{1-\frac{1}{2}, j-3/2}^{n+\frac{1}{2}}}{2\Delta z(j)} + \frac{r_{1+\frac{1}{2}} u_{1+\frac{1}{2}, j-\frac{1}{2}}^{n+\frac{1}{2}} - r_{1-3/2} u_{1-3/2, j-\frac{1}{2}}^{n+\frac{1}{2}}}{\Delta(r_1)^2} \right]$$

Defining the velocity on the right-hand side of the momentum equations at time $n+\frac{1}{2}$ results in

i IS THE RIGHT BOUNDARY
AND j IS THE TOP BOUNDARY OF THE CELL

$$X(i) = \sum_{i=1}^i \Delta X(i)$$

$$Y(j) = \sum_{j=1}^j \Delta Y(j)$$



THE AREA OF CELL (i, j) IN THE i
DIRECTION = $2 \pi X(i) \Delta Y(j)$

THE AREA OF CELL (i, j) IN THE j
DIRECTION = $\pi (X_{(i)}^2 - X_{(i-1)}^2)$

K IS DEFINED AS $(j-1) i \text{ MAX} + i + 1$, AND IS THE INDEX
FOR CELL QUANTITIES

Fig. 2

$$\frac{\rho_{i-\frac{1}{2},j-\frac{1}{2}}}{\Delta t} [(u_{i-\frac{1}{2},j-\frac{1}{2}}^{n+1})^2 - (u_{i-\frac{1}{2},j-\frac{1}{2}}^n)^2] = 2r_{i-\frac{1}{2}}(u_{i-\frac{1}{2},j-\frac{1}{2}}^{n+\frac{1}{2}}) \frac{[P_{i-3/2,j-\frac{1}{2}}^n - P_{i+1/2,j-\frac{1}{2}}^n]}{\Delta(r_i^2)}$$

and

$$\frac{\rho_{i-\frac{1}{2},j-\frac{1}{2}}}{\Delta t} [(v_{i-\frac{1}{2},j-\frac{1}{2}}^{n+1})^2 - (v_{i-\frac{1}{2},j-\frac{1}{2}}^n)^2] = (v_{i-\frac{1}{2},j-\frac{1}{2}}^{n+\frac{1}{2}}) \frac{[P_{i-3/2,i-\frac{1}{2}}^n - P_{j+1/2,i-\frac{1}{2}}^n]}{\Delta z_j}$$

and

$$\begin{aligned} & \frac{\rho_{i-\frac{1}{2},j-\frac{1}{2}}}{\Delta t} [I_{i-\frac{1}{2},j-\frac{1}{2}}^{n+1} - I_{i-\frac{1}{2},j-\frac{1}{2}}^n] \\ &= -P_{i-\frac{1}{2},j-\frac{1}{2}}^n \left[\frac{v_{i-\frac{1}{2},j+\frac{1}{2}}^{n+\frac{1}{2}} - v_{i-\frac{1}{2},j-\frac{3}{2}}^{n+\frac{1}{2}}}{2\Delta z_j} + \frac{r_{i+\frac{1}{2}} u_{i+\frac{1}{2},j-\frac{1}{2}}^{n+\frac{1}{2}} - r_{i-3/2} u_{i-3/2,j-\frac{1}{2}}^{n+\frac{1}{2}}}{\Delta(r_i^2)} \right] \end{aligned}$$

Defining:

$$u_{i-\frac{1}{2},j-\frac{1}{2}}^{n+\frac{1}{2}} = \frac{u_{i-\frac{1}{2},j-\frac{1}{2}}^{n+1} + u_{i-\frac{1}{2},j-\frac{1}{2}}^n}{2}$$

$$v_{i-\frac{1}{2},j-\frac{1}{2}}^{n+\frac{1}{2}} = \frac{v_{i-\frac{1}{2},j-\frac{1}{2}}^{n+1} + v_{i-\frac{1}{2},j-\frac{1}{2}}^n}{2}$$

and

$$P_i^n = \frac{P_{i+\frac{1}{2}}^n + P_{i-\frac{1}{2}}^n}{2}$$

$$P_j^n = \frac{P_{j+\frac{1}{2}}^n + P_{j-\frac{1}{2}}^n}{2}$$

Equation (2), the radial momentum equation, becomes

$$u_{(k)}^{n+1} - u_{(k)}^n = 2\Delta t r_{i-\frac{1}{2}} \pi DY(j) \frac{(PL^n - PRR^n)}{AMX(k)}$$

where

$$PL^n = \frac{P_{(k)}^n + P_{(\text{cell to the left})}^n}{2}$$

$$PRR^n = \frac{P_{(k)}^n + P_{(\text{cell to the right})}^n}{2}$$

Equation (3), the axial momentum equation, becomes

$$v_{(k)}^{n+1} - v_{(k)}^n = \Delta t \pi (r_i^2 - r_{i-1}^2) \frac{(P_{BLO}^n - P_{ABOVE}^n)}{AMX(k)}$$

where

$$P_{BLO}^n = \frac{P_{(k)}^n + P_{(cell\ below)}^n}{2.}$$

$$P_{ABOVE}^n = \frac{P_{(k)}^n + P_{(cell\ above)}^n}{2.}$$

The energy equation (4) becomes

$$I_{(k)}^{n+1} - I_{(k)}^n = \frac{P_{(k)}^n \pi \Delta t}{AMX(k)} \left[\left(\frac{VBLO^n + VBLO^{n+1} - VABOVE^n - VABOVE^{n+1}}{2.} \right) x \right. \\ \left. (r_i^2 - r_{i-1}^2) + [DY(j)] (UL_j^n + UL_j^{n+1} - URR^n - URR^{n+1}) \right]$$

where

$$VBLO = \frac{v_{(k)} + v_{(cell\ below)}}{2.}$$

$$VABOVE = \frac{v_{(k)} + v_{(cell\ above)}}{2.}$$

$$UL = \frac{u_{(k)}^{RC} + u_{(cell\ on\ the\ left)}^{RL}}{2.}$$

$$URR = \frac{u_{(k)}^{RC} + u_{(cell\ to\ the\ right)}^{RR}}{2.}$$

where

$$RC = r_{(i-\frac{1}{2})}$$

$$RR = r_{(i+\frac{1}{2})}$$

$$RL = r_{(i-3/2)}$$

The above equations conserve energy exactly, despite finite difference approximations. However, an adjustment at transmissive boundaries of the grid is necessary. This is a work term, which is also taken into account in ETH (the total energy of the system.) For the transmissive boundaries, the pressure gradient is zero and the velocity at the boundary interface is set equal to the velocity of the cell adjacent to the boundary.

The term subtracted from ETH for the boundary at the right is

$$\frac{P_{(k)} + P_{(\text{cell to the left})}}{2} u_{(k)} r_{i-\frac{1}{2}} \pi \Delta t D Y_{(j)}$$

and the top is

$$\frac{P_{(k)} + P_{(\text{cell below})}}{2} v_{(k)} \pi (r_i^2 - r_{i-1}^2) \Delta t (.5)$$

and the bottom, if transmissive, is

$$\frac{P_{(k)} + P_{(\text{cell above})}}{2} v_{(k)} \pi (r_i^2 - r_{i-1}^2) \Delta t (.5)$$

and is added to ETH. K (in the above equations) refers to the border cell.

The velocity terms in the energy equation for those cells at the transmissive boundaries are, at the right = $u_{(k)} r_{i-\frac{1}{2}}$ and the top = $v_{(k)}$.

Rewriting Eq. (1), the mass transport equation in finite difference form results in

$$\frac{\rho_{(k)}^{n+1} - \rho_{(k)}^n}{\Delta t} = \left[\frac{r_{i-1} \rho_{i-1} u_{i-1}}{r_{i-\frac{1}{2}} \Delta r_i} - \frac{r_i \rho_i u_i}{r_{i-\frac{1}{2}} \Delta r_i} + \frac{\rho_{j-1} v_{j-1} - \rho_j v_j}{\Delta z(j)} \right] \quad (6)$$

where

$$\Delta z(j) = \frac{V_{(k)}}{A_j^z} = \frac{V_{(k)}}{A_{j-1}^z}$$

where A^z for all $j = \pi (r_{(i)}^2 - r_{(i-1)}^2)$,

and

$$V_{(k)} = \text{volume of cell } k = 2\pi r_{i-\frac{1}{2}} \Delta r_i \Delta z_{(j)} \quad (7)$$

multiply both sides of Eq. (7) by r_i results in

$$V_{(k)} r_i = 2\pi r_i \Delta z_j r_{i-\frac{1}{2}} \Delta r_i$$

or

$$V_{(k)} r_i = A_i^r r_{i-\frac{1}{2}} \Delta r_i \quad (8)$$

where A^r = area in the direction perpendicular to the Z axis.

And similarly, multiplying Eq. (7) by r_{i-1} results in

$$V_{(k)} r_{i-1} = 2\pi r_{i-1} \Delta z_j r_{i-\frac{1}{2}} \Delta r_i$$

or

$$V_{(k)} r_{i-1} = A_{i-1}^r r_{i-\frac{1}{2}} \Delta r_i \quad (9)$$

Solving Eqs. (8) and (9) for $r_{i-\frac{1}{2}} \Delta r_i$ and substituting their values into Eq. (6) results in

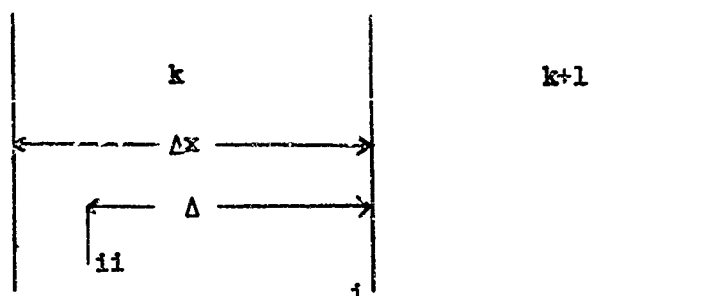
$$\frac{\rho_{(k)}^{n+1} - \rho_{(k)}^n}{\Delta t} = \frac{1}{V_{(k)}} (A_{j-1}^z \rho_{j-1} v_{j-1} - A_j^z \rho_j v_j + A_{i-1}^r \rho_{i-1} u_{i-1} - A_i^r \rho_i u_i)$$

or rewriting in terms of mass,

$$AMX_{(k)}^{n+1} - AMX_{(k)}^n = \Delta t [(Av)_B^z \rho - (Av)_T^z \rho + (Au)_L^r \rho - (Au)_R^r \rho]$$

where AMX = mass of cell k in grams and (Av) = area times a velocity (the velocity is a function of the velocity of the two cells in question.) B refers to the bottom, T to the top, L to the left, and R to the right of cell k . The ρ used is the ρ from the cell from which the flux is coming.

Various techniques for velocity weighting in the mass flux have been tried. Results from these trials are presented in Figs. 5 and 6. The scheme presently being used is as follows. Take the r direction as an example:



The mass to move across i is between i and ii where $\Delta = i - ii$; thus $\Delta = \tilde{u}\Delta t$ where \tilde{u} is the weighted velocity at Δ . Using the first two terms of the Taylor series at a distance of $-\Delta$ from i , we expand

$$u_{(i)} = \frac{u_{(k)} + u_{(k+1)}}{2.}$$

or

$$\tilde{u} = \frac{u_{(k)} + u_{(k+1)}}{2.} + (-\Delta) \frac{(u_{(k+1)} - u_{(k)})}{\Delta x}$$

or

$$\frac{\Delta}{\Delta t} = \tilde{u} = \frac{\frac{u_{(k)} + u_{(k+1)}}{2.}}{\frac{(u_{(k+1)} - u_{(k)})\Delta t}{\Delta x} + 1}$$

if $\tilde{u} > 0$, use $\rho_{(k)}$; if $\tilde{u} < 0$, use $\rho_{(k+1)}$ in the mass flux calculation.

Mass, both components of momentum, and the energy across all four sides of the cell are calculated. By conserving both axial and radial momentum and the total energy, the new velocities are calculated and the new internal energy is then the difference between the total and the kinetic.

A look ahead, two cells in both directions, is done to remove preferential mass transport because of the initial choice of indexing in the r direction first. Take the example where the flux out of the top and right are such that their sum would remove more than the mass in the cell. The code would then assign new fluxes such that the top flux would be its fraction of the total flux out times the mass of the cell, and the right flux would be its fraction of the total flux out times the mass of the cell.

To treat a free surface in the continuous Eulerian scheme, we have chosen to use a density cutoff to limit the mass from flowing through N zones in N time steps. If the mass flow across the free surface results in a density which is less than an input number ($\sim 10^{-3} \rho_0$) the flux is held back. To cut the small precursor ahead of the shock front, the velocities are checked against 10^{-8} cm/sh; if they are smaller than this, they are set to zero.

To ensure that the bottom cells in the projectile will empty as the projectile moves up, a scheme using the ρ and v from the cell above is used to calculate the flux. This is continued until the initial velocity of the bottom cell of the projectile begins to change because of the shock. After this point is reached, no special procedure is used for the bottom cells of the projectile.

Boundary Conditions

These cells adjacent to the axis of symmetry ($r = 0$) have the following boundary conditions; the pressure on the left side of the cell is equal to the pressure of the first cell, and the velocity at the left is set to zero. The pressure at the right interface of a cell whose right neighbor is void is zero, and the velocity is that of the occupied cell; similarly for the case of a void cell above.

The pressure and velocity at a transmissive boundary are the following. The pressure at the boundary is set equal to the pressure at the left or bottom boundary respectively for a right and top transmissive boundary (no acceleration of the border cells) while the velocity is set equal to the border cell velocity.

The top and right boundary of the grid is transmissive; the bottom boundary may be either transmissive or reflective, where the same boundary conditions then will exist as for the top or right and the axis of symmetry.

Two passes are used to solve the change in internal energy due to the work terms. The first pass through, one calculates the new velocities at time $(n+1)$ and simultaneously, the velocities at the interfaces at time (n) . The time (n) interface velocities are also used to evaluate the internal energy contributions due to terms involving these velocities (see Eq. 4.)

In a second pass the time (n+1) interface velocities are calculated and the associated contributions to the internal energies are computed. (A look-ahead feature of two cells in both directions would enable one to use only one pass.)

An option exists for correcting negative internal energies if they arise in phase 1. The cell where the maximum negative internal energy occurred is recorded; assuming the rate of change of internal energy with time is essentially constant, we calculate a smaller time step, such that the new internal energy will be positive. We complete the entire cycle, to time (n+1), now set the time step negative, integrate backward to time (n), and now forward with the new smaller Δt to a revised time (n+1).

Time Control for Code

The time control for the continuous Eulerian is the same as for the particle in cell scheme, with the exception of the r direction. In the z direction:

$$\begin{aligned} \Delta m_z &= \rho \bar{V} A \Delta t & \text{Assume } \bar{V} &= v(k) \\ & & \rho &= \rho(k) \\ \Delta m_z &= AMx(k) \end{aligned}$$

Then

$$\begin{aligned} AMx(k) &= \frac{AMx(k)}{\pi[r_{(i)}^2 - r_{(i-1)}^2]DY(j)} v(k) \pi[r_{(i)}^2 - r_{(i-1)}^2] \Delta t \\ &= AMx(k) v(k) \frac{\Delta t}{DY(j)} \end{aligned}$$

or $|v(k)| \Delta t \leq DY(j)$ such that the flux in the z direction will not empty the cell.

In the r direction, the stability is as follows.

$$\begin{aligned} \Delta m_R &= \rho \bar{u} A \Delta t & \text{Assume } \Delta r_R &= AMx(k) \\ & & \bar{u} &= u(k) \\ & & \rho &= \rho(k) \end{aligned}$$

Then

$$\begin{aligned} \Delta Mx(k) &= \frac{AMx(k)}{2\pi(r_{i-\frac{1}{2}})\Delta r(i)DY(j)} u(k)2\pi r(i)DY(j)\Delta t \\ &= AMx(k)u(k) \frac{r(i)}{r_{i-\frac{1}{2}}\Delta r(i)} \Delta t \end{aligned}$$

or

$$u(k)\Delta t \leq \frac{\Delta r(i)r_{i-\frac{1}{2}}}{r(i)} \leq \frac{TAU(i)}{2\pi r(i)}$$

C = speed of sound, defined as $\sqrt{\gamma P/\rho}$ for a polytropic form of the equation of state and $(\partial P/\partial \rho)^{\frac{1}{2}}_s$ for a real equation of state. Provisions exist for calling either one.

The time control (Δt) conditions are the Courant condition and that the maximum $[|\frac{u}{\Delta x}| \text{ and } |\frac{v}{\Delta y}|] < \frac{1}{\Delta t}$.

Three options for time control are:

1. Code will control the Δt , calculated from the Courant and particle velocity scheme, but at a fraction of stability.
2. The Δt loaded at $t = 0$ will remain constant, provided option for integrating backward in time to remove negative internal energies from phase 1 is not operating.
3. Code will control the Δt , decreasing Δt if

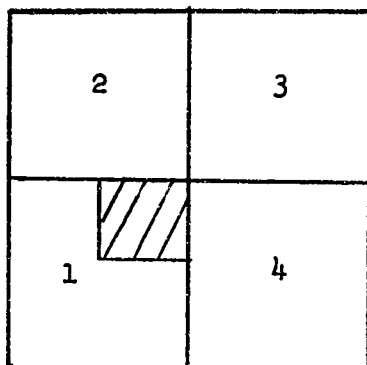
$$\left| \frac{u\Delta t}{\Delta x} \right| \quad \text{or} \quad \left| \frac{v\Delta t}{\Delta y} \right|$$

exceeds FFA (an input number), and increasing Δt if it is less than FFB (an input number.)

The stability check is omitted for a cell if the density of that cell is less than some input number. This prevents isolated debris of high velocity and small masses from controlling the time step.

Corner Coupling

The question investigated here is the correctness of the mass transport which is done neglecting corner coupling. Below is an example of the comparison with a PIC-like transport:



Assume $u = v$ for all four cells and

$$v\Delta t = \frac{\Delta x \text{ or } \Delta y}{2}$$

Where particles in the PIC scheme located in the shaded area will cross into zone 3 in one time step, the OIL code requires two time steps for mass to move into zone 3, first by the path of zone 1 to 2 or 4, and finally to zone 3.

In the case of very small time steps, it is seen that the above approximation is unimportant. We have chosen to run most of our problems at .5 stability. Further, from early test runs of an impact calculation, results did not change appreciably as the factor was varied from one-fourth to one-half to nine-tenths of stability.

The PIC Transport

The changes required to change from a continuous mass transport to the particle transport are:

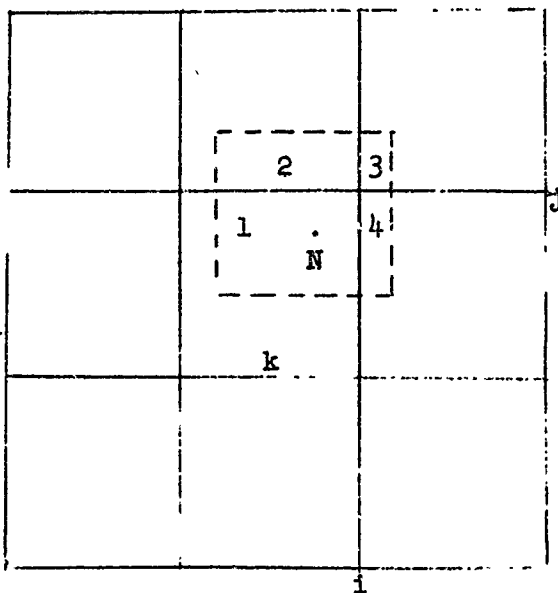
1. The problem number must be negative.
2. The transport and rezone subroutines must be replaced with those for PIC transport.

Following is a brief discussion of the particle (PIC) transport characterizing SHELL. Two scratch tapes are required for the particles. SHELL reads in particle records from one tape, processes the particles updating the coordinates, writes them out on the other tape, and interchanges tape numbers.

Five variables are associated with each particle; the mass (AM), the r coordinate (XL), and Z coordinate (YL), the i-coordinate of the cell where the particle lies (iwl), and the j-coordinate of the cell where the particle

lies (1w2). Thus, one computes the cell number where the particle is by $k = (j-1)*iMAX + i + 1$.

The particles are moved with an area weighted velocity, which is basically a cell placed with the particle at the center. The overlay of this cell on the four cells in question times the velocity of that cell is summed for the four cells, and then the weighted velocity is calculated by dividing through by the total area.



Example: For particle N the u component of velocity used to move the particle is

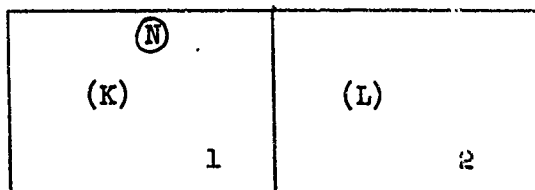
$$\bar{u} = \frac{\sum_{i=1}^4 A_i u_i}{\sum_{i=1}^4 A_i}$$

and the v component is

$$\bar{v} = \frac{\sum_{i=1}^4 A_i v_i}{\sum_{i=1}^4 A_i}$$

The particle is then moved with the area weighted velocities. If the particle does not leave the cell (k), no additional calculations are needed; process the next particle. If the particle leaves cell (k), it carries with it a mass, momentum, and internal energy.

By conserving momentum and total energy, one then changes the quantities in the new cell. No changes except removing the particle mass from the mass of cell (k) are necessary for updating cell (k).



Example: Particle N moves from cell (k) to cell (L). Conserving both axial and radial momentum:

$$(\bar{M}_2)\bar{u}_2 = M_2 u_2 + m u_{11}$$

and

$$(\bar{M}_2)\bar{v}_2 = M_2 v_2 + m v_{11},$$

where the line above signifies the updated variable. Thus,

$$\bar{u}_2 = \frac{M_2 u_2}{\bar{M}_2} + \frac{m u_{11}}{\bar{M}_2}$$

$$\bar{v}_2 = \frac{M_2 v_2}{\bar{M}_2} + \frac{m v_{11}}{\bar{M}_2}$$

However, M_2 is not available at this stage (since \bar{M}_2 has replaced M_2); substitute $M_2 = \bar{M}_2 - m$ results in $\bar{u}_2 = \frac{m}{\bar{M}_2} (u_{11} - u_2) + u_2$; similarly for the v component. Note that u_{11} and v_{11} are set equal to u_1 and v_1 unless there has been an elastic bounce off a reflective boundary (requiring the velocities to change sign to conserve momentum); then u_{11} and v_{11} are set to $-u_1$ and $-v_1$.

To calculate the new internal energy in cell (2) requires that we conserve total energy and momentum, resulting in the expression:

$$M_1 I_1 + \frac{1}{2} M_1 (u_1^2 + v_1^2) + M_2 I_2 + \frac{1}{2} M_2 (u_2^2 + v_2^2) =$$

$$(M_1 - m) I_1 + \frac{1}{2} (M_1 - m) (u_1^2 + v_1^2) + (M_2 + m) \bar{I}_2 + \frac{1}{2} (M_2 + m) (\bar{u}_2^2 + \bar{v}_2^2)$$

Solving for $\bar{I}_2 \bar{M}_2 =$

$$m I_1 + \frac{1}{2} m (u_1^2 + v_1^2) + M_2 I_2 + \frac{1}{2} M_2 (u_2^2 + v_2^2) - \frac{1}{2} \bar{M}_2 (\bar{u}_2^2 + \bar{v}_2^2)$$

Substituting the new values for \bar{u}_2 and \bar{v}_2 from conserving momentum results in

$$\bar{I}_2 = I_2 + \frac{m}{\bar{M}_2} \left\{ I_1 - I_2 + \left[\frac{(u_{11} - u_2)^2 + (v_{11} - v_2)^2}{2} \right] \left(1 - \frac{m}{\bar{M}_2} \right) \right\}$$

Thus, after each particle is moved, the two cells involved are updated. After a particle record has been processed, the particles, with their new coordinates and i and j value of the new cell location, are then written on another tape, which will become the starting conditions for the next transport cycle. An option exists to call rezone if particles leave the top or the right of the grid.

Viscosity

The movement of mass across the cell boundaries give rise to force which is effective in reducing fluctuations that arise from the differencing technique^(ref 1). This is of the form of a "true" viscosity, being proportional to the velocity gradient. This viscosity is present at all times, both in compressions and rarefactions. No additional (that is, a controllable) artificial viscosity is present in this version of oil.

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3.2. Special Subroutines

The REZONE routine for the particle transport is different from that required for the continuous mass transport version. If material leaves the grid out of the right or top, a trigger is set to call REZONE.

The REZONE multiplies all dimensions by two, so that four old cells become one in the new grid. The total number of cells remains the same, and the target is doubled in depth and width by adding mass at the sides and back surface. This scheme does conserve total energy, and by conserving momentum, also, new velocities and internal energy are calculated. This will cause the total internal energy of the system to rise slightly.

For the particle_rezone, it is not necessary to multiply all dimensions by two, but rather change the Δx 's and Δy 's by any prescribed amount. One adds new material with the same density, internal energy, and velocity distributions as are available in CLAM. The number of particles per cell to add is also an input number. For a more complete description, see Reference 3.

A subroutine SETUP is available to generate the initial grid (bypasses the generator code CLAM) if both the target and projectile are of the same density. The projectile must be a right circular cylinder. This routine assumes that all Δx 's are the same and all Δy 's are the same. An asterisk before the symbol implies it is floating point.

SYMBOL	LOCATION	DESCRIPTION
* Z(111)	111	Initial density g/cm^3
* Z(112)	112	Initial pellet velocity cm/sh.
* DX(1)	7845	Δx in cm
* DY(1)	7897	Δy in cm
iMAX	33	Maximum number of zones in the r direction
jMAX	35	Maximum number of zones in the z direction
i1	47	The i value of one radius of the projectile +2
i2	48	The j value of the top of projectile +2
* PROB	1	Any positive number for the problem
* PK(3)	237	Must be a positive number

* PK(4)	238	Set = 1
* PK(5)	239	Right boundary (i) of projectile
* PK(6)	240	Bottom (j)+1 of projectile
* PK(7)	241	Top (j) of projectile
* PK(8)	242	Set = 1
* PK(9)	243	Right (i) boundary of target
* PK(10)	244	Bottom (j)+1 of target
* PK(11)	245	Top (j) of target

And the usual input data to start OIL code from a CLAM tape.

A. Example of Input for OIL Using the Subroutine SETUP.

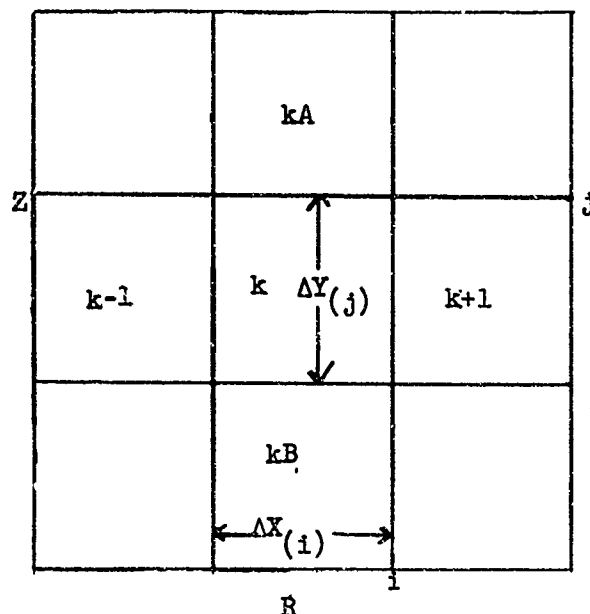
(OIL INPUT)

GAS		GAMMA= 1.5		RHO TARGET =		RHO PELLET =1.						
1	23531.		0.		0.							
	11211.01											
	1111.											
	11111.											
	7511.	-3										
	13821.	-2.5										
	23871.	6.		9.		20.		1.		24.		21.
	245132.											
2	33124.											
2	35132.											
	784511.											
	789711.											
	7711.											
	271-1.											
	821-1.											
	651-1.											
	7111.											
	1421.	.5										
	2421.	-50.										
	8621.5	1.5										
	147120.											
2	4728.	22.										
	11311.	-2										
1	3510.	4.		4.		16.		16.				
1	7211.	44										

This will make a binary tape that is equivalent to that made by the example in CLAM (see Section 2.1). See Section 5.2 for format for CARDS subroutine.

3.3. Logic of OIL

The logic involved in following a given cell (K) from time t to t+Δt, or from cycle n to n+1.



$$k = (j-1)iMAX + i + 1$$

$$TAU(i) = \pi(x_{(i)}^2 - x_{(i-1)}^2)$$

Fig. 3

1. CDT Routine

Here one calculates the pressure (P) for cell (k) where $P(k) = f[\rho(k), I(k)]$ where $\rho(k) = \frac{AMZ(k)}{TAU(i)DY(j)}$. The speed of sound $(\gamma P/\rho)^{\frac{1}{2}}$ or $(\partial P/\partial \rho)^{\frac{1}{2}}$ is then calculated and the Courant condition for stability, that $\frac{C\Delta t}{\min(\Delta x \text{ or } \Delta y)} < \frac{1}{2(\gamma_{\max}-1)}$ and the particle velocity criteria that the $\max\left[\left|\frac{u}{\Delta x}\right| \text{ and } \left|\frac{v}{\Delta y}\right|\right] < \frac{1}{\Delta t}$ are calculated. From these stability checks a new Δt is calculated or Δt remains the same (see options under description of common for OIL, symbol CABIN.) The cycle number and the time are now advanced. (A radiation time step is also calculated, although radiation is not being used in this version of the code.)

2. EDIT Routine

The OIL code has four separate editing-like routines all included in the routine called EDIT. A section called short print displays the time, cycle number, the total internal and kinetic energy and total mass, and various other integral quantities such as momenta, and mass in various

angles (see Sect. 3.4.) A plot routine is also available in the EDIT routine which places an (x) (in an equal cell size grid, corresponding to the actual grid in OIL) if there is any mass in the cell; thus it is a display of mass movement from time to time. A long print routine may also be called for that edits on each page a column from the OIL grid which contains the coordinates of the column, and the cell quantities as a function of the row coordinates (see column identification in Sect. 5.2 FORTRAN listing of EDIT.)

The last option is a dump routine which dumps all necessary data for restarting the problem. This data may also be used for the automatic plotting routines.

Various input numbers (see Sect. 3.4) specify the frequency that these routines will be called for, and an input number specifying a cycle number to stop.

In the following discussion, please refer to Fig. 3.

3. PHL Routine

Here we integrate the two momentum equations and the change in internal energy due to the work terms. No material is moved at this time, and the transport terms are dropped. Using the new pressures and the time step which were computed in CDT, we now prepare to integrate the equations.

PL(j), the pressure at interface (i-1) and uL(j), the velocity at interface (i-1) are available from the previous column sweep on i-1.

$$PL(j) = \frac{P(k) + P(k-1)}{2}$$

$$uL(j) = \frac{r_{i-3/2} u_{(k-1)}^n + r_{i-1/2} u_{(k)}^n}{2}$$

The PBLO term, which was the PABOVE for cell (kB) and VBLO, VABOVE for cell kB, is also available for interface j-1.

$$PBLO = \frac{P(k) + P(kB)}{2}$$

$$VBLO = \frac{v(k) + v(kB)}{2}$$

Now we calculate terms at interface i and j:

$$\begin{aligned} PR &= \frac{P(k+1) + P(k)}{2.} & URR &= \frac{r_{i-\frac{1}{2}} u(k)^n + r_{i+\frac{1}{2}} u(k+1)^n}{2.} \\ P_{ABOVE} &= \frac{P(k) + P(kA)}{2.} & VABOVE &= \frac{v(k)^n + v(kA)^n}{2.} \end{aligned}$$

Now we can integrate the two momentum equations

$$\rho \frac{\partial u}{\partial t} = - \frac{\partial P}{\partial r}$$

or

$$u_{(k)}^{n+1} = u_{(k)}^n + \left(\frac{PL(j)^n - PRR^n}{AMX(k)} \right) 2r_{i-\frac{1}{2}} \pi \Delta t DY(j)$$

and

$$\rho \frac{\partial v}{\partial t} = - \frac{\partial P}{\partial z}$$

or

$$v_{(k)}^{n+1} = v_{(k)}^n + \left(\frac{PBLO^n - PABOVE^n}{AMX(k)} \right) \pi (r_i^2 - r_{i-1}^2) \Delta t$$

Now one can add the work term due to velocities at cycle N to the change in internal energy.

$$\begin{aligned} I_{(k)}^{n+1} &= I_{(k)}^n + \frac{P(k)}{AMX(k)} [(uL_{(j)}^n - uRR^n) \pi \Delta t DY(j) \\ &\quad + \left(\frac{VBLO^n - VABOVE^n}{2.} \right) \pi (r_i^2 - r_{i-1}^2) \Delta t] \end{aligned}$$

$$\rho \frac{\partial I}{\partial t} = - P \left(\frac{\partial v}{\partial z} + \frac{1}{r} \frac{\partial ur}{\partial r} \right)$$

Now we make one more pass through the grid, this time omitting the momentum equations but calculating the velocity terms at the interface, where again we only have to calculate the data at interface i and j:

$$\begin{aligned}
 uL(j) &= \frac{r_{i-3/2} u_{(k-1)}^{n+1} + r_{i-1/2} u_{(k)}^{n+1}}{2} & VABOVE &= \frac{v_{(k)}^{n+1} + v_{(kA)}^{n+1}}{2} \\
 uRR &= \frac{r_{i-1/2} u_{(k)}^{n+1} + r_{i+1/2} u_{(k+1)}^{n+1}}{2} & VBLO &= \frac{v_{(k)}^{n+1} + v_{(kB)}^{n+1}}{2}
 \end{aligned}$$

and then add to $I_{(k)}^{n+1/2}$ the work terms due to velocities at cycle $n+1$.

$$\begin{aligned}
 I_{(k)}^{n+1} &= I_{(k)}^{n+1/2} + \frac{P_{(k)}^n}{AMX(k)} [(uL_{(j)}^{n+1} - uRR^{n+1}) \pi \Delta t DY(j) \\
 &\quad + (\frac{VBLO^{n+1} - VABOVE^{n+1}}{2}) \pi (r_1^2 - r_{i-1}^2) \Delta t]
 \end{aligned}$$

The specific internal energy is checked during both passes for negative values. If a negative value is found, we assume dI/dt is constant over the time step, and recompute a new Δt (not placing it in the Δt storage) that will prevent I from going negative. After the completion of integrating all values to cycle $n+1$, an option exists for removing these negative energies. If one selects the option, Δt is set equal to $-\Delta t$, the code integrates backward two passes to cycle n , then replacing Δt with the new, smaller Δt and now forward in time with two passes using the smaller Δt .

4. PH2 Routine

Here we move the mass and approximate the transport terms in the momentum and energy equations that were omitted in Phase 1.

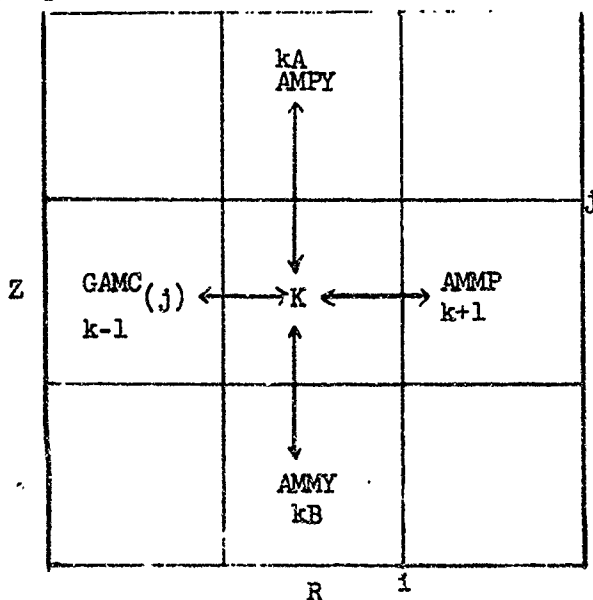


Fig. 4

AMMP = mass flow across the right boundary
 AMUR = radial momenta across the right boundary
 AMVR = axial momenta across the right boundary
 DELER = total specific energy across the right boundary
 AMBY = mass flow across the bottom boundary
 AMBU = radial momenta across the bottom boundary
 AMBV = axial momenta across the bottom boundary
 DELEB = total specific energy across the bottom boundary
 GAMC(j) = mass flow across the left boundary
 FLEFT(j) = radial momenta across the left boundary
 YAMC(j) = axial momenta across the left boundary
 SIGC(j) = total specific energy across the left boundary
 AMPY = mass flow across the top boundary
 AMUT = radial momenta across the top boundary
 AMVT = axial momenta across the top boundary
 DELET = total specific energy across the top boundary

Again, following a typical cell k (Fig. 4) the masses, the momentas, and the energies are now available at the left and bottom boundaries of cell (k) from the previous column sweep and the cell below.

Now, begin by calculating the mass flow at the top of cell (k).

$$V_{\text{ABOVE}} = \frac{v(k) + v(kA)}{2}$$

then form

$$\frac{V_{\text{ABOVE}}}{1 + \left[\frac{v(kA) - v(k)}{\Delta y(j)} \right] \Delta t}$$

as the weighted velocity to use in the flux equation and store it in V_{ABOVE} . If it is positive, use $\rho(k)$; if it is negative, we use $\rho(kA)$. Now calculate the mass flow across the top as $\Delta m_T = \rho(M) A(j) V \Delta t$ where

$$\begin{aligned} \bar{V} &= V_{\text{ABOVE}} \\ M &= \text{donor cell} \end{aligned}$$

or

$$\Delta m_T = \Delta m_F = \frac{\Delta m_x(m)}{\Delta y(j)} V_{\text{ABOVE}}(\Delta t)$$

Now we calculate the mass flow at the right boundary of cell k. u_{RR} is defined as $\frac{u(k) + u(k+1)}{2}$ then form

$$\frac{u_{RR}}{1 + \left[\frac{u(k+1) - u(k)}{\Delta x(i)} \right] \Delta t}$$

as the weighted velocity to use in the flux equation and store it in u_{RR} . The mass flow across i is then

$$\Delta m_R = \rho(M) A(i) \bar{u} \Delta t$$

where

$$\bar{u} = u_{RR}$$

$$N = i \text{ value of donor cell}$$

$$M = \text{donor cell}$$

$$PIDTS = \frac{1}{\pi \Delta t}$$

or

$$\Delta m_R = \Delta m_P = \frac{\Delta m_x(M)}{\tau(N)} \frac{x(i)}{PIDTS} 2(u_{RR})$$

Now check to see if these masses will more than empty the cell, since it is possible that the left and bottom flux were both negative. Search cells ahead in both directions to remove preferential mass movement. As an example, suppose the flux at the TOP (Δm_F) is positive and the flux at the right is positive, where their sum is larger than the mass in cell (k); then normalize the fluxes the following way. The flux out of the top is

$$\frac{F_T[\Delta m_x(k)]}{F_T + F_R}$$

and the flux out of the right is equal to

$$\frac{F_R[\Delta m_x(k)]}{F_T + F_R}$$

where F is a symbol for the mass flux.

The momenta associated with these masses are now computed. The sign of the flux specifies the zone where the mass came from, thus at the top; the radial component equals $AMUT = AMPY(u(N))$ and the axial $AMVT = AMPY(v(N))$ where (N) = cell number of the donor cell.

The momenta for the right is $AMUR = AMMP(u(N))$ for the radial and $AMVR = AMMP(v(N))$ for the axial component, where again (N) = cell number of the donor cell.

The total specific energy that those mass fluxes carry is also calculated at this time; for the top it is equal to

$$DELET = I_{(n)} + \frac{u_{(N)}^2 + v_{(N)}^2}{2}$$

and for the right it is equal to

$$DELER = I_{(n)} + \frac{u_{(N)}^2 + v_{(N)}^2}{2}$$

where again (N) = cell number of the donor cell.

The mass now in cell (k) is equal to $DEIM = AMX(k) + GAMC(j) + AMMY - AMPY - AMMP$ which equals the original mass plus the mass flow across the left, the bottom, and less the mass flow across the top and the right.

The total axial momenta that have come into or left cell (k) is = $SIGMV = YAMC(j) + AMMV - AMVT - AMVR$ = the momenta crossing the left boundary plus the momenta crossing the bottom boundary less the momenta crossing the top and the right boundary.

The total radial momenta that has come into or left cell (k) is $SIGMU = FLEFT(j) + AMMU - AMUT - AMUR$ = momenta crossing the left and bottom boundary less the momenta crossing the top and the right.

Similarly we calculate the total energy that these fluxes have carried = $DELEK = (GAMC(j)) SIGC(j) + (AMMY) Deleb - (AMPY) Delet - (AMMP) Deler$ = the mass times the total specific energy at the left plus the similar term for the bottom less the similar terms for the top and the right.

Now by conserving momenta and total energy, calculate the new specific internal energy and velocities of cell (k).

$$MU_L + MU_B - MU_T - MU_R + MU_k = (\text{TOTAL MASS})\bar{u}$$

where total mass = DEIM and

$$Mv_L + Mv_B - Mv_T - Mv_R + Mv_k = (\text{TOTAL MASS})\bar{v}$$

and the new specific internal energy

$$I_{(k)} = \frac{E_L + E_B - E_T - E_R + E_k}{\text{DEIM}} - \frac{\bar{u}^2 + \bar{v}^2}{2}$$

and now AMX(k) set = to DEIM.

The subscripts L, B, T, R, refer to the left, bottom, top, and right. Now the values that we calculated at the right for cell (k) are now set to the left values for cell (k+1) and the top values for cell (k) now become the bottom values for cell (kA).

The limits of the DO loops on i and j are i1 and i2. A check is done to see if mass or energy has moved beyond i1 or i2 and then the counter is increased by 1. This check is also done in PH1. By using this scheme, we process only the active mesh at all times, in all the subroutines.

3.4. List of Common (OIL)

The location refers to the location of that symbol relative to the beginning of common. Since the beginning of common is assigned the same location for each subroutine, a program (CARDS) is available for changing any word in common. The z block is first in common for OIL.

Note that if one should change the dimensions of the arrays, he must be careful and also make the necessary changes of the locations in the following tables.

<u>Symbol</u>	<u>Location</u>	<u>No. of Words</u>	<u>Units</u>	<u>Description</u>
AID	706	1	--	Not used, since this is a one-material (x) code
AIX	707	3500	jerks/g	Specific internal energy (x) for cell (k)
AM	4207	130	g	Mass of particle (N) for PIC transport only
AMD	4337	1	--	Not used in this one-material code
AMK UR(16)	220	15	g	Storage (EDIT) for summing masses in given angles for editing
AMX	4338	3500	g	Total (x) mass in cell k
AREA	7838	1		Tag, used in PH2 (EUL and PIC)
BIG	7839	1	--	Not used
BOUNCE	7840	1	--	Tag used in particle PH2
CABLN	82	1	--	If < 0 code controls Δt but at Z(139) of instability

Caution: You must load a Δt for this option

If = 0, code will control the Δt , decreasing Δt if $\left| \frac{u\Delta t}{\Delta x} \right|$ or $\left| \frac{v\Delta t}{\Delta y} \right|$ exceed FFA (an input number) and increasing Δt if $\left| \frac{u\Delta t}{\Delta x} \right|$ or $\left| \frac{v\Delta t}{\Delta y} \right|$ is less than FFB (an input number)

This holds provided SN \neq 0

If > 0, DT loaded will remain constant.

DDXN	7841	1	--	Not used
DDVK	7842	1	--	Not used
DKE	7843	1	--	Not used
DVK	7844	1	--	Not used
DX	7845	52	cm	$DX(i) = X(i) - X(i-1)$
DY	7897	100	cm	$DY(j) = Y(j) - Y(j-1)$
E	7997	1	--	Not used
FD	7998	1	--	Not used
FS	7999	1	--	Not used PH2
FX	8000	1	--	Not used
OUT	8001	1	--	Tag in particle PH2
P	8002	3500	jerks/cm ³	Material pressure in cell k
PABOVE	11502	1	jerks/cm ³	$\frac{P_{(k)} + P_{(cell\ above)}}{2}$

PBLO	11503	1	jerks/cm ³	$\frac{P_{(k)} + P_{(\text{cell below})}}{2.}$
PIDTS	11504	1	1/cm.sh	$= \frac{1.}{\Delta t \pi DY(j)} \text{ in (PH1);}$ $\frac{1}{\pi \Delta t} \text{ in PH2}$
PK UR(31)	235	15	g.cm/sh	Radial momentum in certain angles for editing
PL,PR,GAMC PR(100),SIG(C)	405	200	Many	Pressure in PH1, flux. in PH2, etc.
PPABOV	11505	1	--	Not used
PRR	11506	1	jerks/cm ³	$\frac{P_{(k)} + P_{(\text{cell to the right})}}{2.}$
FUL	11507	1	--	Not used
QDT	11508	1	--	Not used
QK	250	15	g.cm/sh	Axial momentum in certain angles for editing
RC	11509	1	cm	$[(X(i) + X(i-1))]/2. \text{ in PH1}$
ReZ	11510	1	--	If > 0 and ReZFct > 0, then PH2 (EUL) will call subroutine REZONE, ReZ set in PH2.
RHO	11511	1	g/cm ³	Density
RL	11512	1	--	Not used
RR	11513	1	cm	$[X(i) + X(i+1)]/2 \text{ in PH1}$
SIG	11514	1	cm	Minimum Δx or Δy in CDT routine
SIGN = QOOOFL	11515	1	0	Not used
SWITCH	11516	1	--	Not used
TAB	205	15	--	Tan(α) (Table of, used in EDIT routine)
TABLM	11517	1	--	Not used
TAU	11518	52	cm ²	$= \pi(X_{(i)}^2 - X_{(i-1)}^2) = \text{area}$
TAUDTS	11570	1	cm ² sh	$= \text{TAU}(\perp) \Delta t \text{ in PH1.}$
TAUDTX	11571	1	--	Not used
U	11572	3500	cm/sh	R component of velocity in cell (k)
UK	15072	1	cm/sh	R component of velocity in cell (k) used in PIC transport

UL,UR, Fleft. etc.	205	200	cm ² /sh	$\frac{U(k)RC + U(k-1)RCC}{2.}$ where $RCC = \frac{X_{(i-1)} + X_{(i-2)}}{2.}$ and $RC = \frac{X_{(i)} + X_{(i-1)}}{2.}$
URR	15073	1	cm ² /sh	[U(k)RC + U(k+1) RR]/2.
UT	15074	1	--	Signal in PH1 (decrease Δt)
UU	15075	1	sh	New Δt if PH1 integrates back for I < 0, set originally to 10 ⁻¹⁵ .
UUU	15076	1	--	Not used
UTEF	15077	1	cm/sh	R velocity component used to move particle when using PIC transport
UVMAX	15078	1	l/sh	Max velocity /Min(ΔX or ΔY)
V	15079	3500	cm/sh	Axial (Z) component of velocity for cell (k)
VABOVE	18579	1	cm/sh	[V(k) + V(cell above)]/2.
VBLO	18580	1	cm/sh	[V(k) + V(cell below)]/2.
VEL	18581	1	--	Used as a tag in PH1 and Max(γ-1) in CDT, and tag in TUL. PH2
VK	18582	1	cm/sh	Axial component of velocity in cell (k) for PIC transport
VT	18583	1	--	Not used
VTEF	18584	1	cm/sh	Z velocity component used to move particle when using PIC transport
VV	18585	1	--	Not used
VVABOV	18586	1	--	Not used
VVBLO	18587	1	--	Not used
W2	18588	1	--	Not used
W3	18589	1	--	Not used
WPS	18590	1	--	Working Storage
ws	18591	1		↓
WSA	18592	1		
WSB	18593	1		
WSC	18594	1		

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XX	152	53	cm	XX(2) = X(1)
XL	18595	130	cm	R coordinate of particle N
XLF	18725	1	--	Fraction of area on left to use in velocity weighting for PIC PH2.
XN	18726	1	cm	R coordinate of particle N at cycle (n-1)
XR	13727	1	--	Fraction of area to the right to use in velocity weighting for PIC PH2.
X1	73	1	--	Not used
X2	79	1	--	Not used
Y	606	100	cm	Y(j) = top dimension of zone (i,j)
YL	18723	130	cm	Z coordinate of particle N
YLW	18358	1	--	Fraction of area below to be used in velocity weighting for PIC PH2.
YN	13359	1	cm	Z coordinate of particle N at cycle (n-1)
YU	18360	1	--	Fraction of area above to be used in velocity weighting for PIC PH2.
YY (YY(2) = X(1))	605	1	cm	YY(j+1) = Y(j)
Y1	80	1	--	Not used
Y2	81	1	--	Not used
Z	1	150	--	See pages where Z(1) through Z(150) are defined
ZMAX	18361	1	cm	An up-to-date value of largest Y coordinate of all particles used in PIC PH2.
i	18362	1	--	<u>Index (Working Storage)</u>
i1	18363	1		
iN	18364	1		
iR	18365	1		
iWS	18366	1		
iWSA	18367	1		
iWSB	18368	1		
iWC	18369	1	--	
i1	47	1	--	The right boundary of the active grid + 2, MAX(i1) = iMAX

12	48	1	--	The top boundary of active grid + 2 MAX(12) = JMAX
13	49	1	--	Not used
14	50	1	--	Not used
1W1	18870	130	--	= (i) of the cell (k) where particle (N) is located, used in PIC PH2.
J	19000	1	--	Index (Working Storage)
JN	19001	1		
JP	19002	1		
JR	19003	1		
K	19004	1	--	Index of cell defined such that $k = (j-1) \text{IMAX} + i+1$
KDT	41	1	--	If KDT = 0, Δt has changed if $\neq 0$ Δt remains constant
KN	19005	1	--	Index (Working Storage)
KP	19006	1		
KR	19007	1		
KPM	19008	1		
L	19009	1		
M	19010	1		
MA	19011	1		
MB	19012	1		
MC	19013	1		
MD	19014	1		
ME	19015	1		
MZ	19016	1	--	Set by input, used in EDIT for length of Z block to write on tape, present value = 150
N	19017	1	--	Index(Working Storage)
NK	19018	1		
NKMAX	19019	1	--	
NK1	19020	1		
NO	19021	1		
NR	19022	1	--	Maximum number of radiation cycles/ hydro cycle calculated in CDT $NR \leq NRM$

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IW2	19023	130	--	= (j) value of cell (k) where particle (N) is located, used in PIC PH2.
IZ	1	150	Many	Fixed value of Z block
Fleft	205	100		Equivalenced, used for flux terms
YAMC	304	100		in Eulerian continuous (PH2)
SIGC	504	100		

<u>Location</u>	<u>Symbol</u>	<u>Units</u>	<u>Description</u>
Z(1)	PROB	--	Problem number (if positive, this is an OIL run; if negative, this is a PIC run)
Z(2)	CYCLE	--	Cycle number (floating point value)
Z(3)	DT	sh	$\Delta t_{\text{hydro}} = t^n - t^{n-1}$
Z(4)	PRINTS	--	Cycle frequency for short print
Z(5)	PRINTL	--	Cycle frequency for long print
Z(6)	DUMPT7	--	Cycle frequency for binary tape dumps
Z(7)	CSTOP	--	Cycle number at which problem stops
Z(8)	PIDY	--	$= \pi = 3.1415927$
Z(9)	TMZ	g	Total (X + .) mass at $t = 0$ (calculated in CLAM)
Z(10)	GAM	--	Not used
Z(11)	GAMD	--	$1./(\gamma_- - 1)$ Computed in Input
Z(12)	GAMX	--	$1./(\gamma_x - 1)$
Z(13)	ETH	jerk	Total energy (computed in CLAM for $t=0$.) Changed in PH1 at transmissive boundaries and in PH2 if mass leaves the system, and by radiation flow out of the system.
Z(14)	FFA	--	Upper limit for stability and to calculate Δt , only if CABLN = 0.
Z(15)	FFB	--	Lower limit for stability and to calculate Δt , only if CABLN = 0.
Z(16)	TMDZ	g	Total (.) mass ($t = 0$) calculated in CLAM
Z(17)	TMXZ	g	Total (x) mass ($t = 0$) calculated in CLAM
Z(18)	XMAX	cm	$= X(iMAX)$
Z(19)	TXMAX	cm	$2(XMAX) t = 0$. calculated in CLAM
Z(20)	TYMAX	cm	$2(YMAX) t = 0$. calculated in CLAM
Z(21)	AMDM	g	Min(.) particle mass/2.; calculated in CLAM

Z(22)	AMXM	g	Min (x) particle mass/2. Calculated in CLAM
Z(23)	DNN	--	$(ETH - E)^{n-NPC}/ETH$
Z(24)	DMIN	--	IE (ECK) Note Z(76) > DMIN, problem will stop and the edit routine will call dump.
Z(25)	FEF	--	Not used
Z(26)	DTNA	sh	Δt^{n-1}
Z(27)	CVIS	--	If < 0, bottom boundary is transmissive; otherwise reflective boundary.
Z(28)	NPR	--	Index (Working storage)
Z(29)	NPR1	--	" "
Z(30)	NC	--	Cycle number in fixed point.
Z(31)	NPC	--	Number of cycles between short prints
Z(32)	NRC	--	Index
Z(33)	IMAX	--	Maximum number of zones in R direction
Z(34)	IMAXA	--	IMAX + 1
Z(35)	JMAX	--	Maximum number of zones in Z direction
Z(36)	JMAXA	--	JMAX + 1
Z(37)	KMAX	--	(IMAX)(JMAX) + 1
Z(38)	KMAXA	--	KMAX + 1
Z(39)	NMAX	--	Total number of particles + 1, generated in CLAM for PIC problem only.
Z(40)	ND	--	Total number of (.) particles + 1 generated in CLAM
Z(41)	KDT	--	Defined previously
Z(42)	ixMAX	--	Not used
Z(43)	NOD	--	Index
Z(44)	NOPR	--	Index
Z(45)	NiMAX	--	New iMAX before adding new zones
Z(46)	NjMAX	--	New jMAX before adding new zones
Z(47)	i1	--	Defined previously
Z(48)	i2	--	Defined previously
Z(49)	i3	--	Not used
Z(50)	i4	--	Not used
Z(51)	N1	--	Scratch tape number for particles if this is a PIC run.
Z(52)	N2	--	Scratch tape number for particles if this is a PIC run.

43

Z(53)	N3	--	Number of particle records generated if this is a PIC run.
Z(54)	N4	--	Number of particles-1 per record (MAX = 127) if this is a PIC run.
Z(55)	N5	--	Not used
Z(56)	N6	--	Number of particles on last particle record if this is a PIC run
Z(57)	N7	--	Not used
Z(58)	N8	--	Not used
Z(59)	N9	--	Not used
Z(60)	N10	--	= i value of zone that is controlling Δt
Z(61)	N11	--	= j value of zone that is controlling Δt
Z(62)	NFM	--	= maximum number of Rad cycles/Hydro (input number)
Z(63)	TRAD	sh	$NR \cdot \Delta t \text{ Rad} = \Delta T \text{ Hydro}$; not used in this version
Z(64)	XNRG	jerks	Total energy of (x) material
Z(65)	SN	--	If = 0 code will decrease Δt to correct for $I < 0$, if $\neq 0$, those $I < 0$ are left alone
Z(66)	DXN	--	Not used
Z(67)	RADER	g-cm/sh	Total positive radial momentum (x only)
Z(68)	RADET	g-cm/sh	Total positive axial momentum (x only)
Z(69)	RADEB	g-cm/sh	Total positive radial momentum (x) for material under target
Z(70)	DTRAD	--	Not used
Z(71)	REZFCT	--	If = 0, PH2 will not trigger rezone
Z(72)	RSTOF	--	Not used in continuous version
Z(73)	SHELL	--	Not used
Z(74)	BBOUND	--	Not used in this version
Z(75)	TOZONE	g/cm ³	Minimum density for mass flow at free surface
Z(76)	ECK	energy check	$\left[\left(\frac{ETH - E}{ETH} \right)^n - \left(\frac{ETH - E}{ETH} \right)^{n-NPC} \right] / NPC$
Z(77)	SBOUND	--	Fraction of Δ in mass weighting velocity EUL PH2 ~ 1.0
Z(78)	X1	--	Not used
Z(79)	X2	--	Not used
Z(80)	Y1	--	Not used

Z(81)	Y2	--	Not used
Z(82)	CABLN	--	Already defined
Z(83)	VISC	--	Not used
Z(84)	T	sh	Total time up to cycle N, $t^n = t^{n-1} + \Delta t$
Z(85)	GMAX	--	Maximum of γ_x or γ .
Z(86)	WSGD	--	γ .
Z(87)	WSGX	--	γ_x and $(\gamma_{\max} - 1)$ in CDT
Z(88)	GMADR	--	$\gamma / (\gamma - 1)$
Z(89)	GMAXR	--	$\gamma_x / (\gamma_x - 1)$
Z(90)	S1	--	Not used
Z(91)	S2	--	Not used
Z(92)	S3	--	Not used
Z(93)	S4	--	Not used
Z(94)	S5	--	Not used
Z(95)	S6	--	Not used
Z(96)	S7	--	Not used
Z(97)	S8	--	Used in CLAM only
Z(98)	S9	--	Not used
Z(99)	S10	--	Not used
Z(100)	f		Mass thrown away (PH2) continuous transport
Z(101)	jerks		Total energy thrown away
Z(102)	g-cm/sh		Total radial momentum thrown away
Z(103)	g-cm/sh		Total axial momentum thrown away
Z(104)	jerks		Energy (internal) added to system when internal is set to 0 if $I < 0$
Z(105)	--		Not used
Z(106)	--		Not used
Z(107)	--		Not used
Z(108)	--		Not used
Z(109)	--		Not used
Z(110)	jerks/g		Critical energy $E(S)$, same value as Z(122)
Z(111)	g/cm ³		Initial density of material
Z(112)	cm/sh		Initial velocity of pellet
Z(113)	--		Epsilonics for emptying pellet $\approx .01$
Z(114)	--		Not used

50

Z(115)	g/cm ³	Density (ρ_0)	For equation of state
Z(116)	--	a	
Z(117)	Jerks/g	E_0	
Z(118)	--	b	
Z(119)	jerks/cm ³	A	
Z(120)	-	V_s	
Z(121)	-	--	
Z(122)	jerks/g	E_s	
Z(123)	--	α	
Z(124)	--	β	
Z(125)	--	--	
Z(126)	jerks/cm ³	B	↓
Z(127)	--	Not used	
Z(128)			
Z(129)			
Z(130)			
Z(131)			
Z(132)			
Z(133)			
Z(134)			
Z(135)			
Z(136)			
Z(137)			
Z(138)	g/cm ³	Density check if $\rho(k) < Z(138)$ stability check for cell (k) is bypassed.	↓
Z(139)	--	Percent of instability, used in CDT if CABIN < 0 $\approx .5$	
Z(140)	--	Not used	
Z(141)	--		
Z(142)			
Z(143)			
Z(144)			
Z(145)			
Z(146)			
Z(147)	--	j (of pellet-target interface) at t = 0	
Z(148)	A	} C(speed of sound = $A + BP^6$ where $A = C_0$ and P is in megabars	
Z(149)	B		
Z(150)	e		

See Ref. 5 for a more detailed description of the equation of state.
For condensed states,

$$P = \left[a + \frac{b}{\frac{E}{E_0 \eta^2} + 1} \right] \frac{E}{V} + A\mu + B\mu^2 ;$$

for expanded states,

$$P = aE\rho + \left[\frac{bE\rho}{\frac{E}{E_0 \eta^2} + 1} + A\mu e^{-\alpha(\frac{V}{V_0} - 1)} \right] e^{-\beta(\frac{V}{V_0} - 1)^2}.$$

$$\rho = \frac{1}{V} \quad \eta = \frac{\rho}{\rho_0} , \text{ and } \mu = \eta - 1 .$$

E = specific internal energy

Condensed form for states where

$$\frac{V}{V_0} \leq f. \quad \text{and} \quad E < E_S$$

expanded form for states where

$$\frac{V}{V_0} > 1 \quad \text{and} \quad E > E_S$$

The following table contains the constants for the above equations.

EQUATION OF STATE DATA

Loc.	Z(115)	Z(116)	Z(117)	Z(118)	Z(119)	Z(120)	Z(121)	Z(122)	Z(123)	Z(124)	Z(125)	Z(126)
Def.	ρ	a	E_0	b	A	V_s	0	E_s	α	β	0	B
Unit	g/cm ³		jerks/ g		jerks/ cm ³			jerks/ g				jerks/ cm ³
W	19.17	.5	2.25 (-5)	1.04	3.08 (-4)	1.11	0.	1.135 (-6)	10.	10.	0.	2.5 (-4)
CU	8.9	.5	3.25 (-5)	1.5	1.39 (-4)	1.13	0.	1.38 (-6)	5.	5.	0.	1.1 (-4)
FE	7.86	.5	9.5 (-6)	1.5	1.279 (-4)	1.21	0.	2.44 (-6)	5.	5.	0.	1.5 (-4)
AL	2.7	.5	5. (-6)	1.63	7.52 (-5)	1.1	0.	3. (-6)	5.	5.	0.	6.5 (-5)
BE	1.845	.55	1.75 (-5)	.62	1.173 (-4)	1.1	0.	1. (-5)	5.	5.	0.	5.5 (-5)
TI	4.51	.5	7. (-6)	.60	1.03 (-4)	1.09	0.	3.5 (-6)	5.	5.	0.	5. (-5)
NI	8.86	.5	9. (-6)	1.33	1.912	1.11	0.	2.85 (-6)	5.	5.	0.	1.5 (-4)
MO	10.2	.5	4.5 (-5)	1.02	2.713 (-4)	1.08	0.	2.3 (-6)	5.	5.	0.	1.65 (-4)
TH	11.63	.4	2.5 (-6)	.36	5.31 (-5)	1.15	0.	2. (-6)	9.	0.88	0.	5. (-5)
CH ₂	.92	.6	7. (-6)	0	7.5 (-6)	1.1	0.	2.4 (-6)	10.	5.	0.	2. (-6)
PB	11.34	.4	1.5 (-6)	2.3	4.664 (-5)	1.1	0.	2.6 (-7)	13.	15.	0.	1.5 (-5)

	Z(148)	Z(149)	Z(150)
	C_0	A	ϵ
AL	3.28	4.4	.45
PB	2.0	2.42	.43
Plastic	2.86	2.2	.45
FE	4.04	3.4	.42
W	4.1	3.4	.42

Where

$$C = C_0 + AP^6$$

 C_0 in 10^5 cm/sec

P in megabars

4. TEST PROBLEMS

4.1. Technique

A series of one-dimensional impact problems was undertaken to select the most appropriate scheme for the velocity weighting in the mass transport equation. An iron projectile traveling at a velocity of 1×10^6 cm/sec striking an iron target was used as the standard test problem. The projectile had five zones of 1-cm each; the target had 95 zones of 1-cm each. The equation of state used was that for iron.⁽⁵⁾

Four different velocity weighting techniques were investigated to determine the best \bar{u} to use in the mass transport expression $\rho \bar{u} \Delta t$. These are illustrated below, for the case where the flow is from cell k to cell $(k+1)$.

The four different approaches were as follows:

1. The donor cell scheme: $\bar{u} = u_{(k)}$

2. The Rich scheme⁽⁶⁾
(no density terms) $\bar{u} = u_{(k)} + \frac{u_{(k+1)} - u_{(k-1)}}{4}$

3. The OIL scheme $\bar{u} = \frac{\frac{u_{(k)} + u_{(k+1)}}{2}}{\left[1 + \frac{[u_{(k+1)} - u_{(k)}] \Delta t}{\Delta x}\right]}$

and 4, another scheme where

$$\bar{u} = \frac{u_{(k)} + u_{(k+1)}}{2} \left[1 - \frac{[u_{(k+1)} - u_{(k)}] \Delta t}{\Delta x}\right]$$

The donor cell, Fig. 5, looks very good in the neighborhood of the shock front; however, behind the front (the rarefaction side) instability sets in. Rich's scheme (without his density weighting) is very similar to scheme 4. The present scheme 3 now used in the OIL code was chosen as the best representation of the shock front and the rarefaction (Fig. 6). However, the results indicate the possibility of using the donor cell (scheme 1) near the shock front and one of the other three schemes in the rest of the problem.

4.2. Comparison of SHELL and OIL

An iron projectile, 3-cm-diam by 3-cm long, strikes an iron target at a velocity of 4×10^6 cm/sec. The problem was run using both SHELL (PIC transport) and OIL (continuous mass transport.) The starting conditions and cell sizes were the same for both. There were 72 cells in the projectile and 2976 in the target, and 16 particles per cell were used in the SHELL run.

Figure 7 is a plot of the total positive axial momentum and the total positive radial momentum (in the target) vs time for the two problems. Figure 8 shows the shock pressure vs position as a function of time for the two problems. The agreement between the two techniques is very good, with the SHELL scheme overshooting the theoretical shock pressure by a somewhat larger amount. Figures 9 and 10 are mass flux plots of the two schemes vs zone number, again at 45° . Figure 11 is a pressure and compression plot of the two schemes as a function of zone number along a ray 45° from the initial center at the projectile target interface. Figure 12 is the same information at a later time of 9.2 μ sec. Throughout all the comparisons, the two schemes are seen to be in very good agreement, with the SHELL scheme exhibiting some undesirable oscillatory behavior due to the discrete nature of the particle population in the cells. Finally, the SHELL run required longer machine time by a factor of 15 over the OIL run.

4.3. One-dimensional Test Problems

Figure 13, a comparison of the two schemes for a plane wave free expansion problem, indicates good agreement between OIL and theory. A hot gas extending to 30 cm with a ρ of $.8 \text{ g/cm}^3$, a γ of $5/3$, and specific internal energy of 9×10^{-3} jerks/g, with a rigid wall on the left, was suddenly released and free to rarefy to zero pressure. The slight rise in mass in the leading edge is due to the constraint adopted for preventing mass transport to diffuse n cells in n time steps (see Section 3.1).

A second "shock tube" problem consisted of a rigid boundary on the left, a hot gas of $\rho = .8 \text{ g/cm}^3$, a γ of $5/3$, and a specific internal energy of 9×10^{-3} jerks/g, which was allowed to shock a cold region extending from 30 cm to 40 cm, consisting of a ρ of 1.2 g/cm^3 and a γ of $5/3$. In

addition to the OIL and SHELL versions of this flow, a third comparison was made using a plate-geometry version of a one-dimensional Lagrangian (SPUTTER) code. The results are given in Figs. 14 and 15, where pressure and velocity are plotted as functions of distances in both the hot and cold material. The comparison between OIL and SPUTTER is very good, while SHELL, again, exhibits an oscillatory behavior.

Figure 16 shows the initial setup of a spherical blast problem which was run using the three codes, SHELL, OIL, and a spherical version of SPUTTER. Figure 17 is the pressure versus position for the three cases, and Fig. 18 is a plot of density versus position. In the latter plot particularly, SPUTTER displays a decided advantage over the two Eulerian methods. This is due to the fact that the SPUTTER (Lagrangian) zoning was on an equal mass basis and the resolution is accordingly better at big R. Hence, there are many zones to represent the density discontinuity or shock location. The SHELL and OIL results are very similar, with the OIL results being substantially smoother.

Figures 19 and 20 are pressures vs position along the r and z axis for the SHELL and OIL codes. The results for OIL are especially interesting in that they are very nearly spherical, where the results for SHELL are not as smooth and are slightly different along the two axes, partially because of the preferential movement of particles that exist in the PIC transport of SHELL.

The spherical character of the OIL solution is also borne out by further tests, such as the close agreement, to about one-tenth of one percent, of the two components of velocity along a ray 45° from the axis of symmetry. The fact that an initially spherical problem remains spherical is gratifying as evidence that the differencing or transport schemes are not introducing significant preferential treatments in the radial or axial directions.

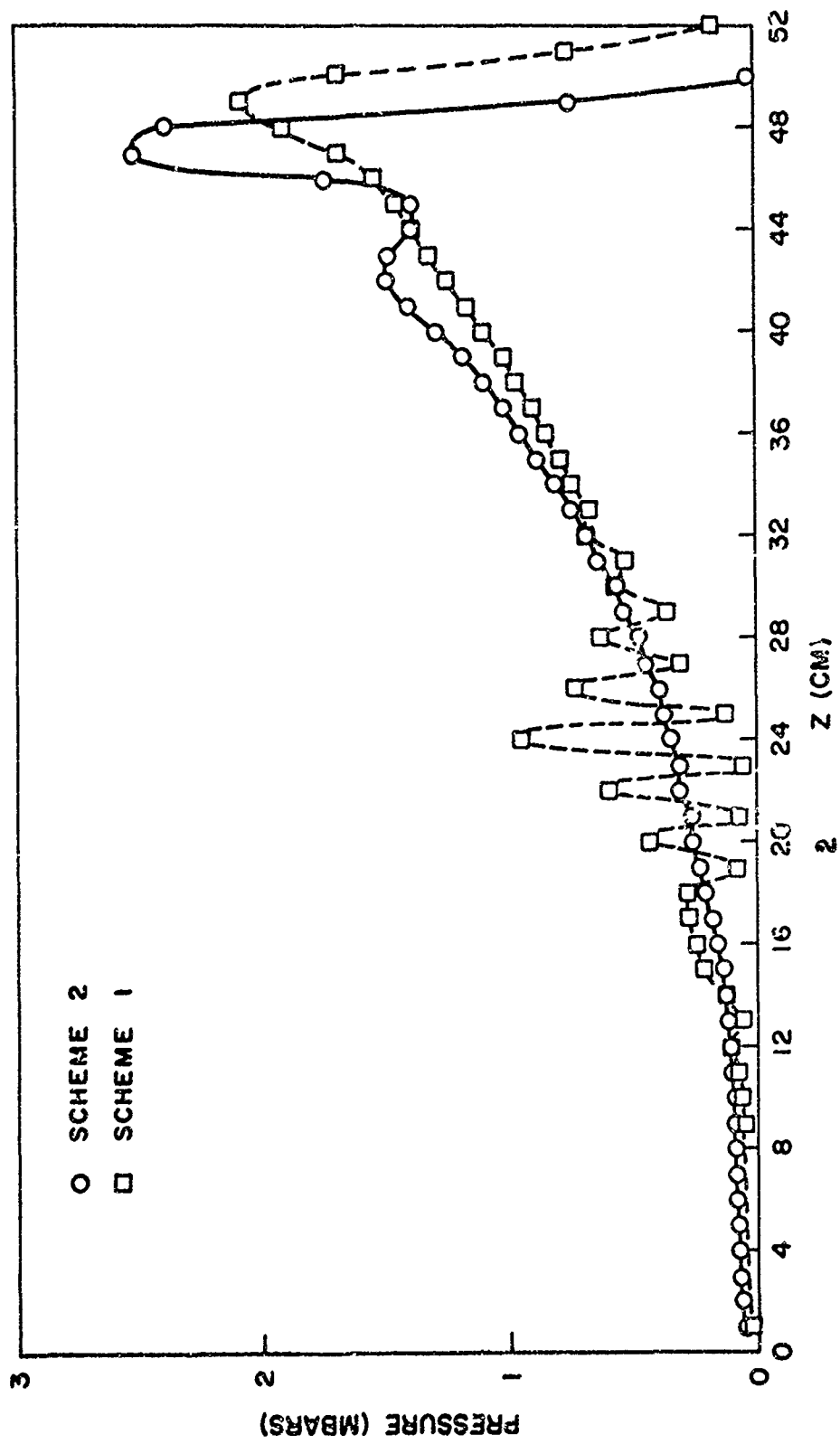


Fig. 5

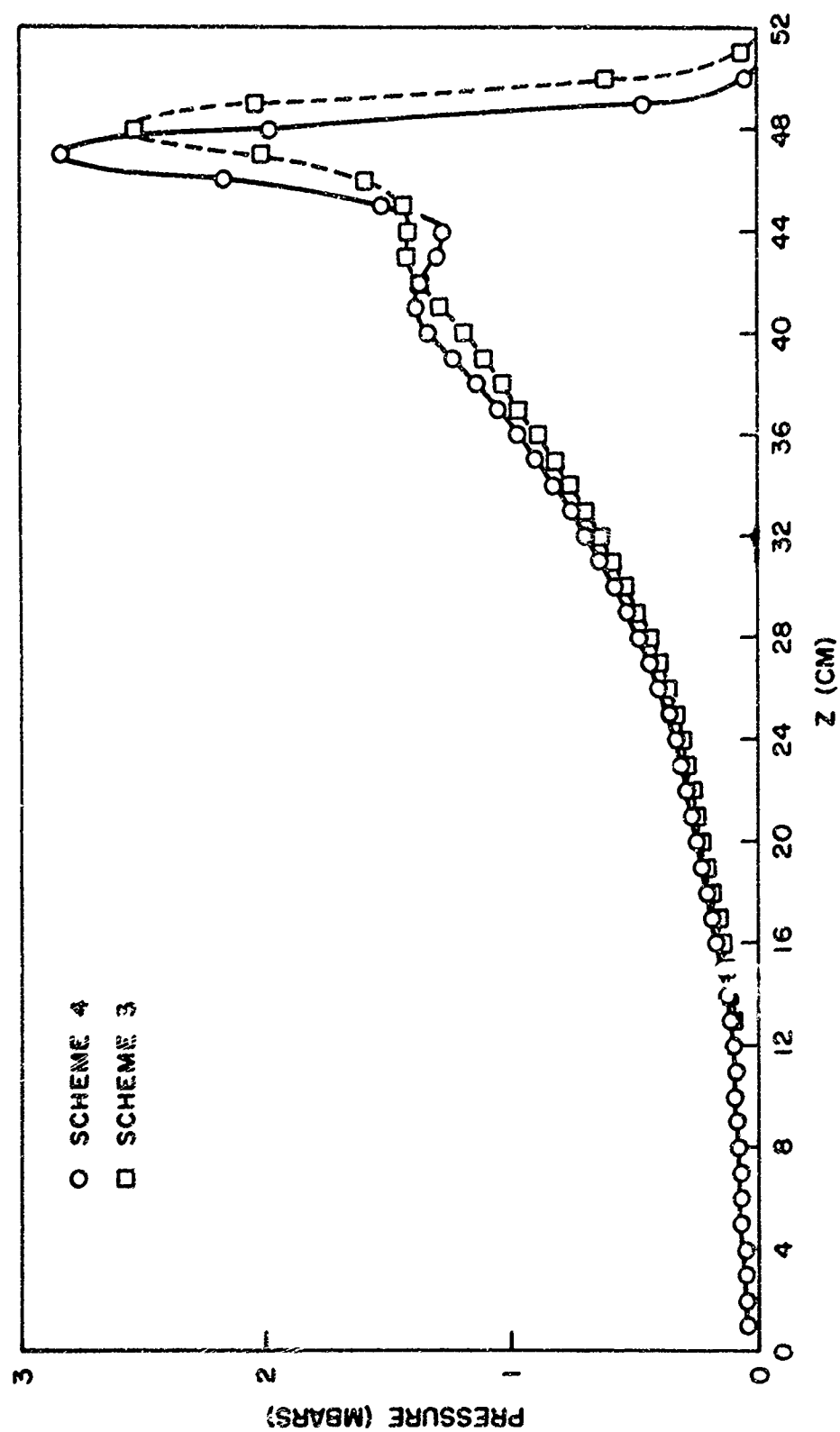


Fig. 6

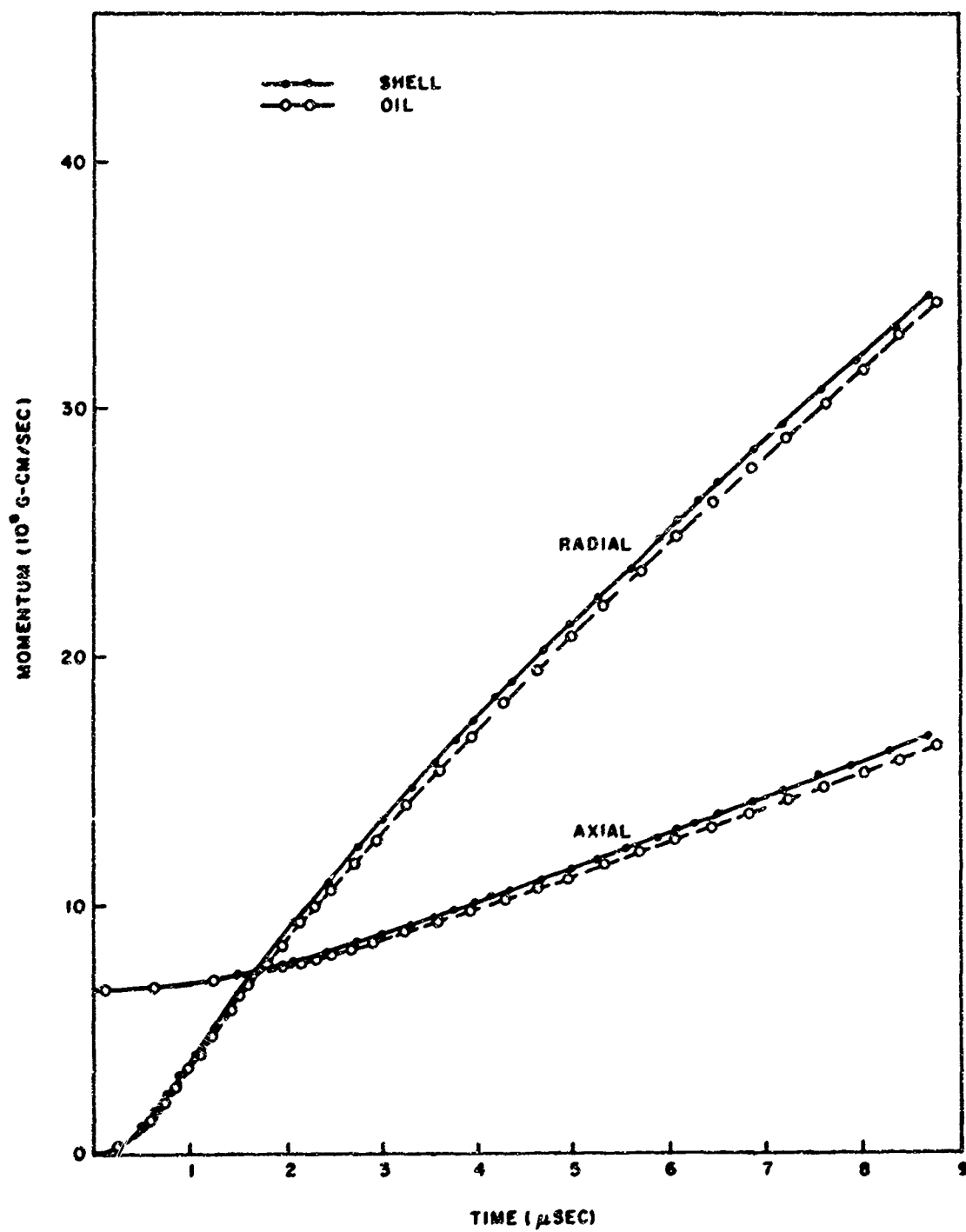


FIG. 7

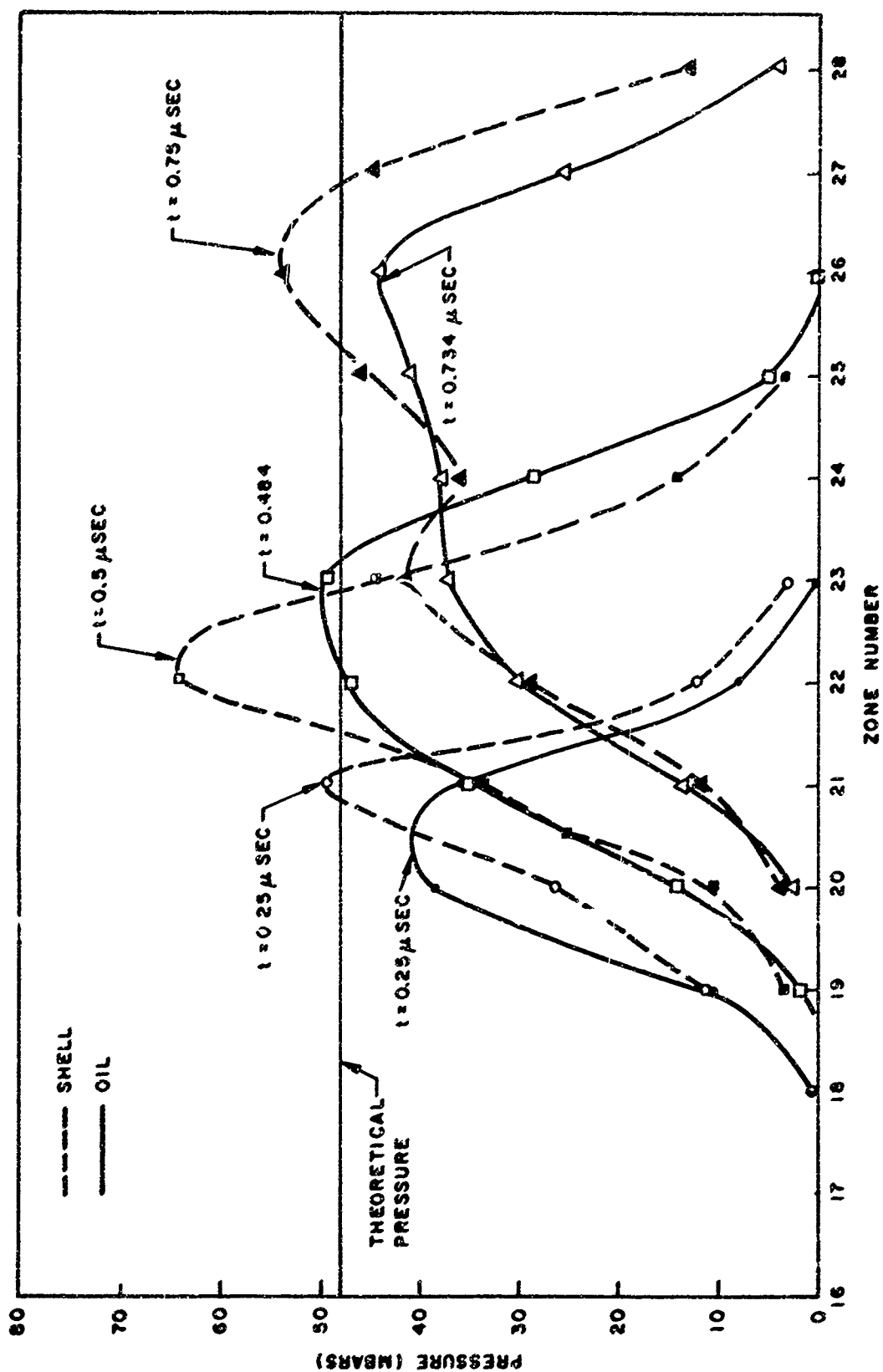


Fig. 8

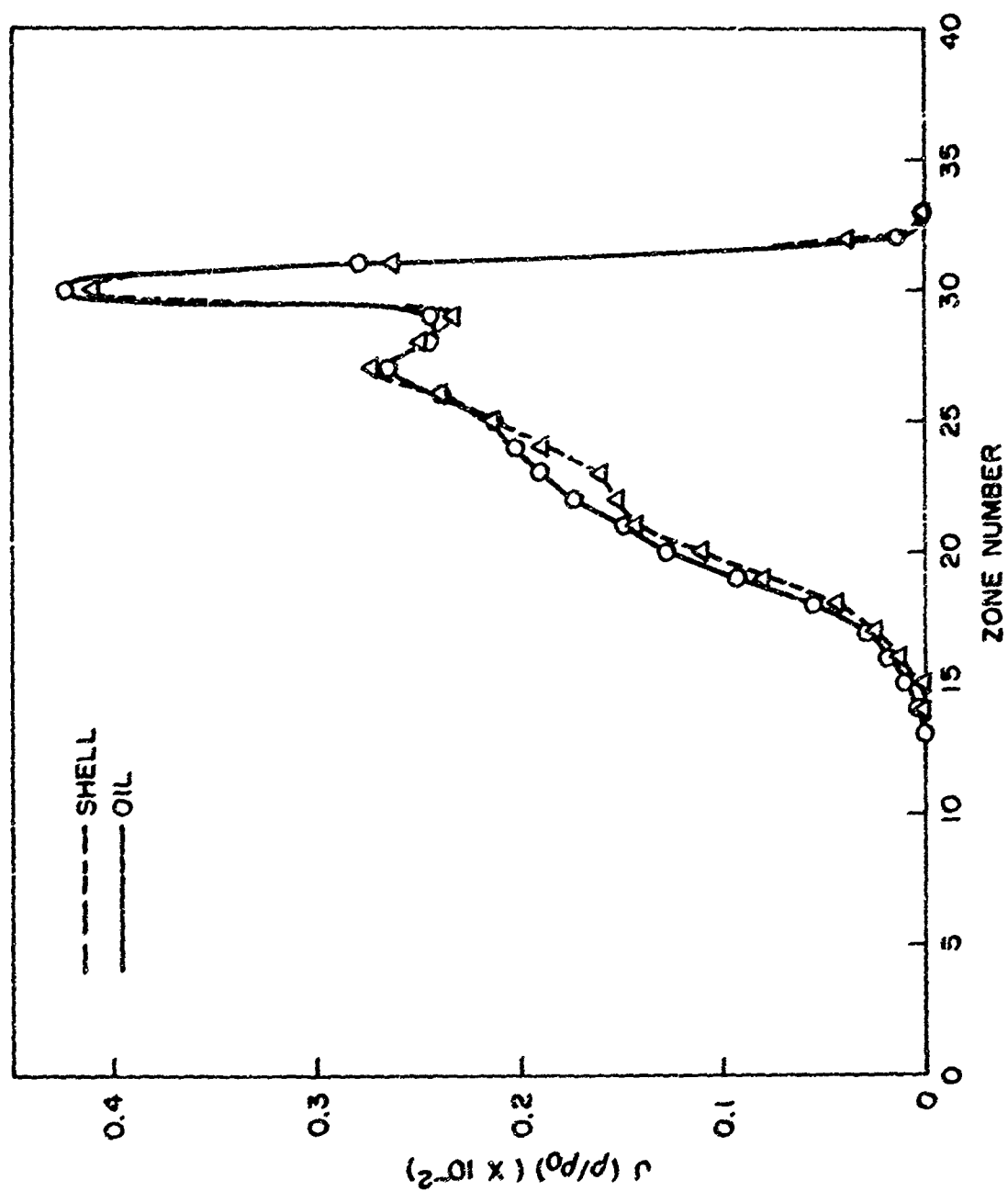


Fig. 9

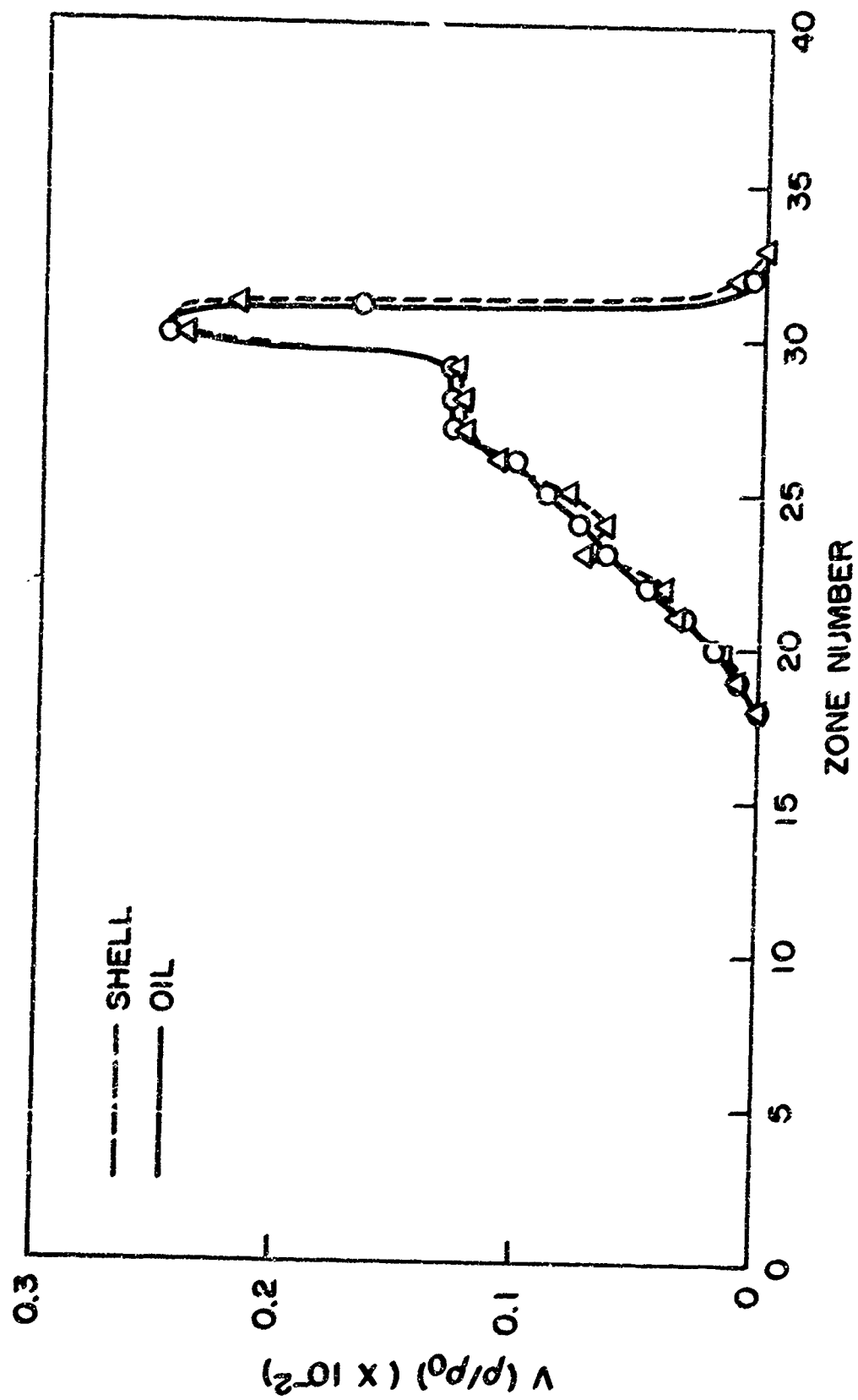


Fig. 10

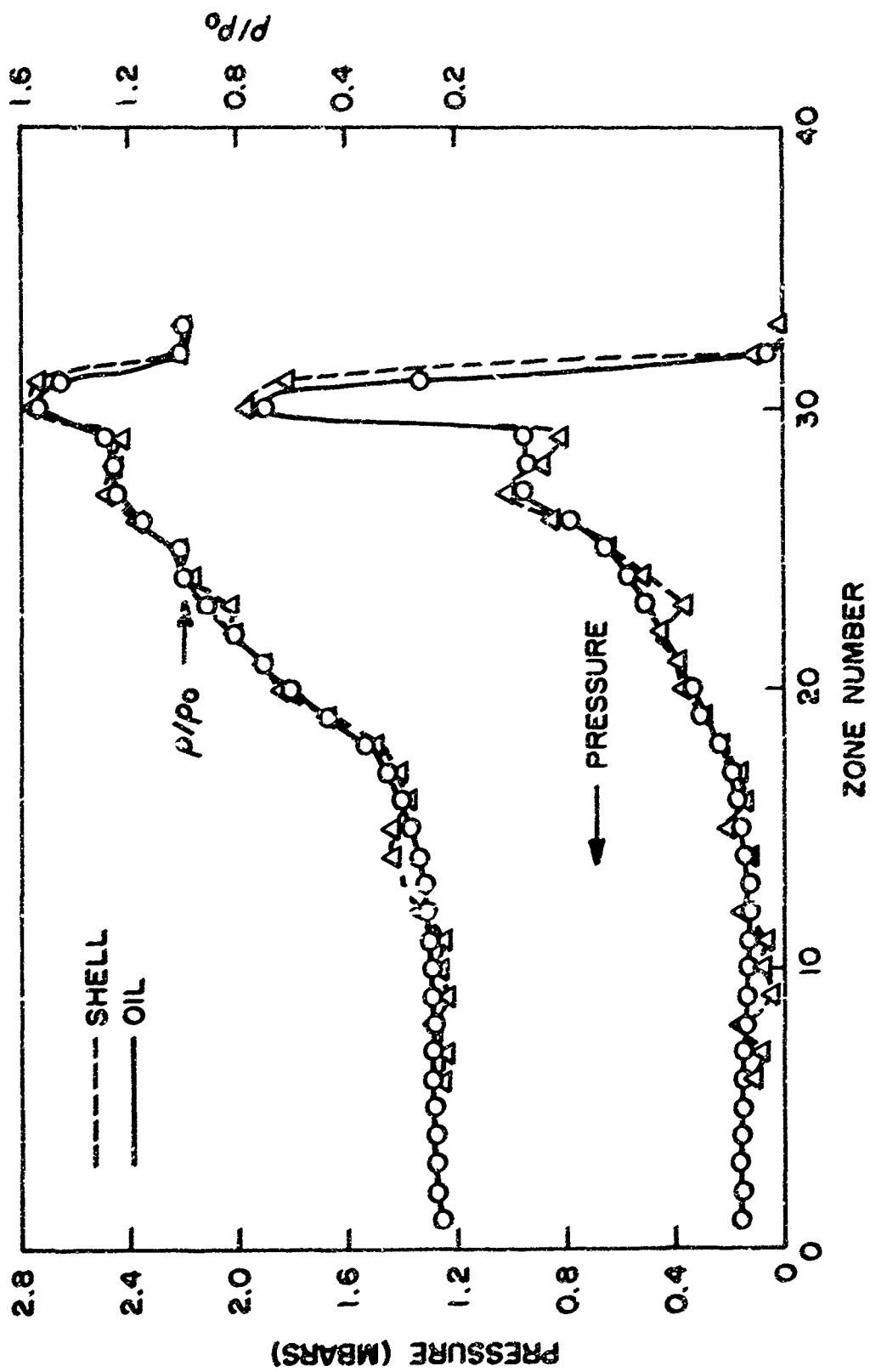


Fig. 1

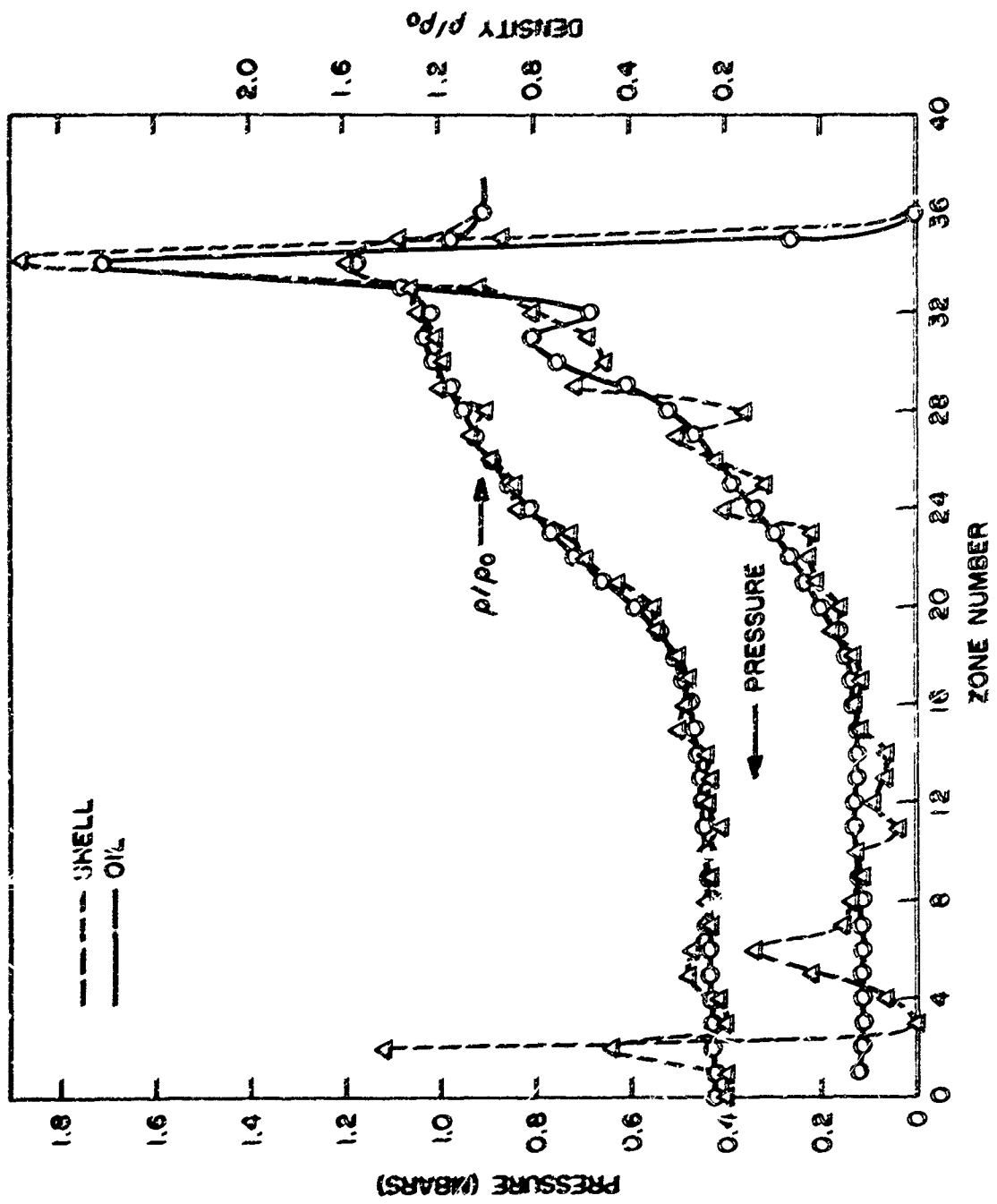


Fig. 12

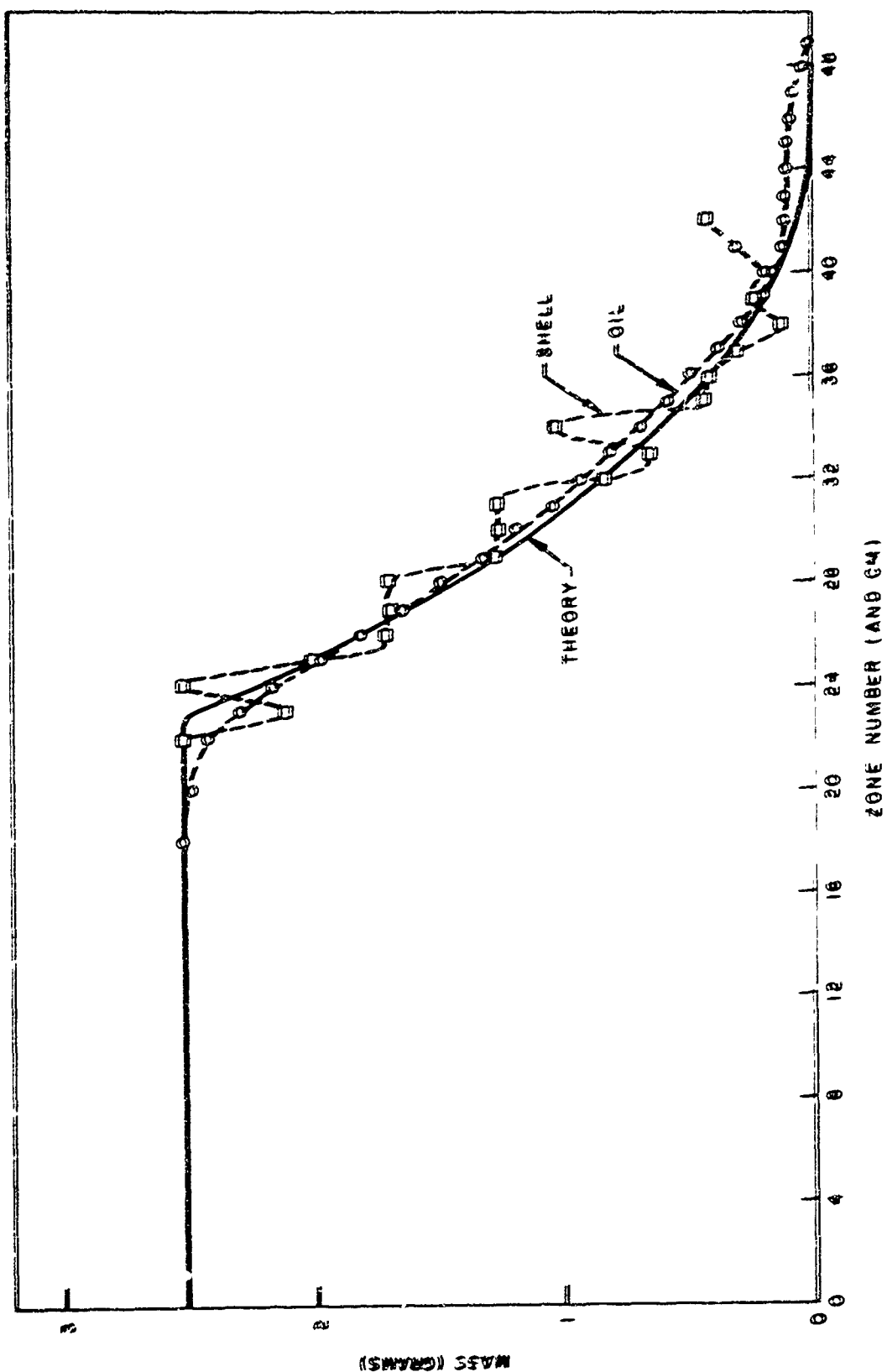
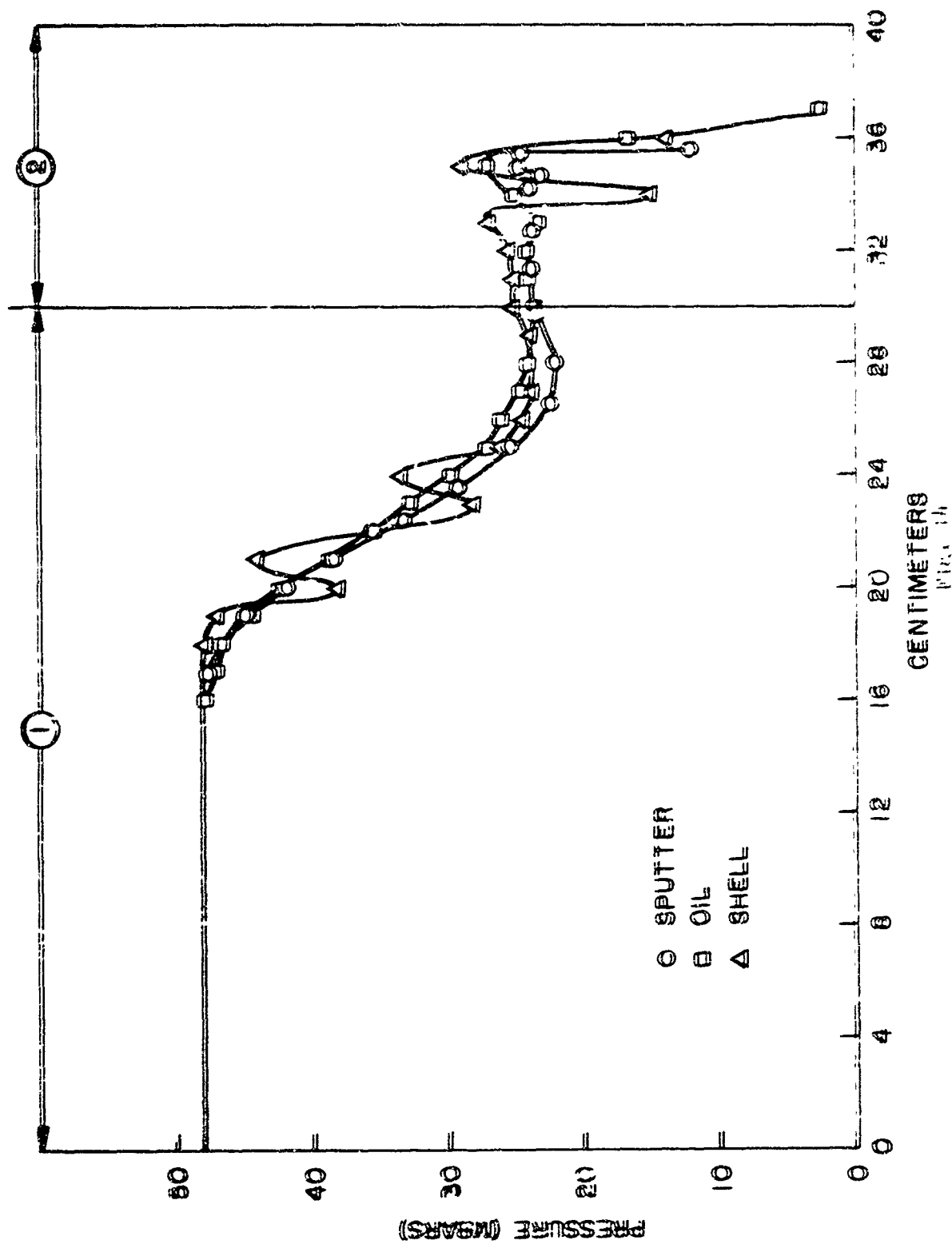


Fig. 12



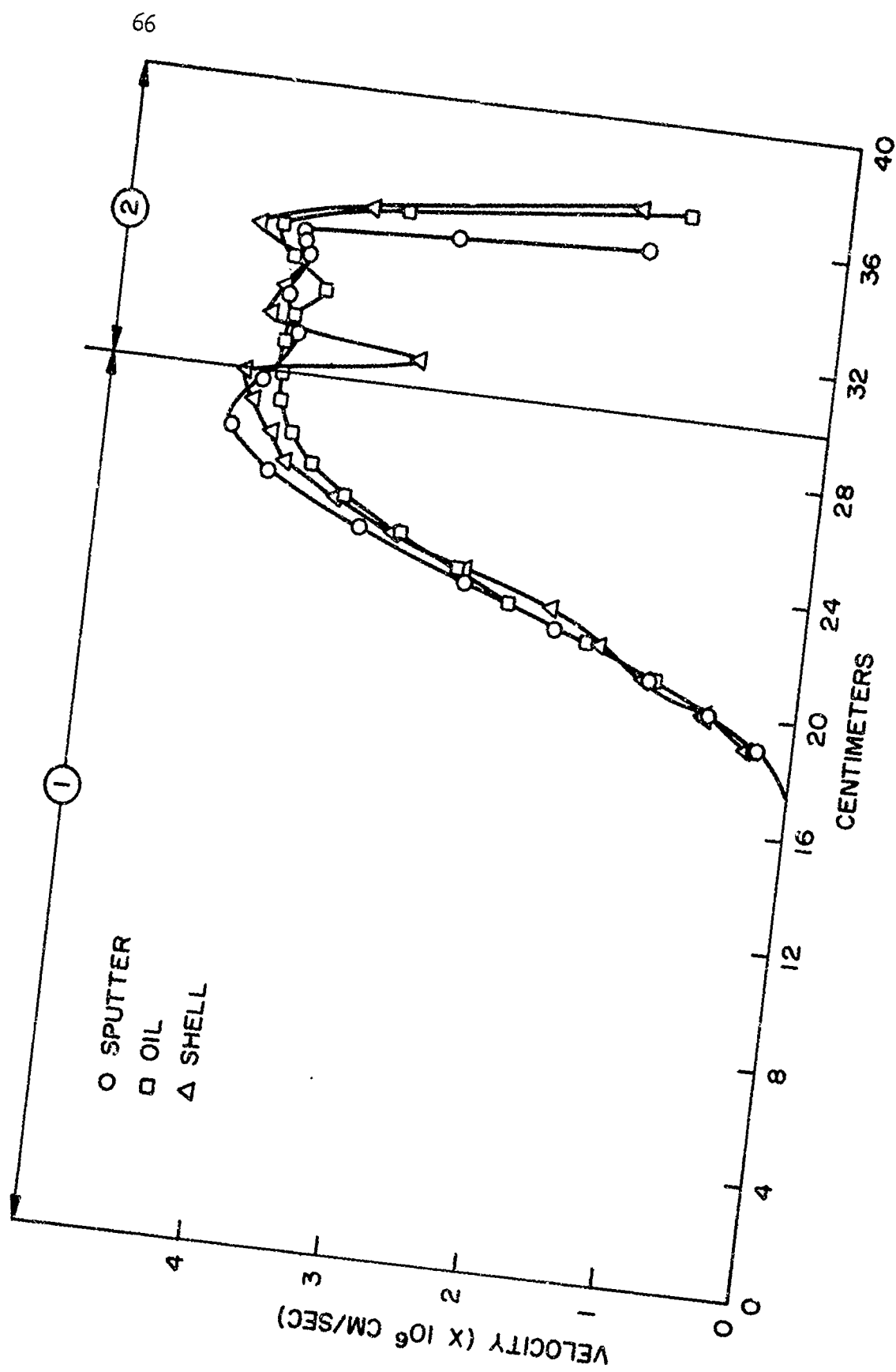
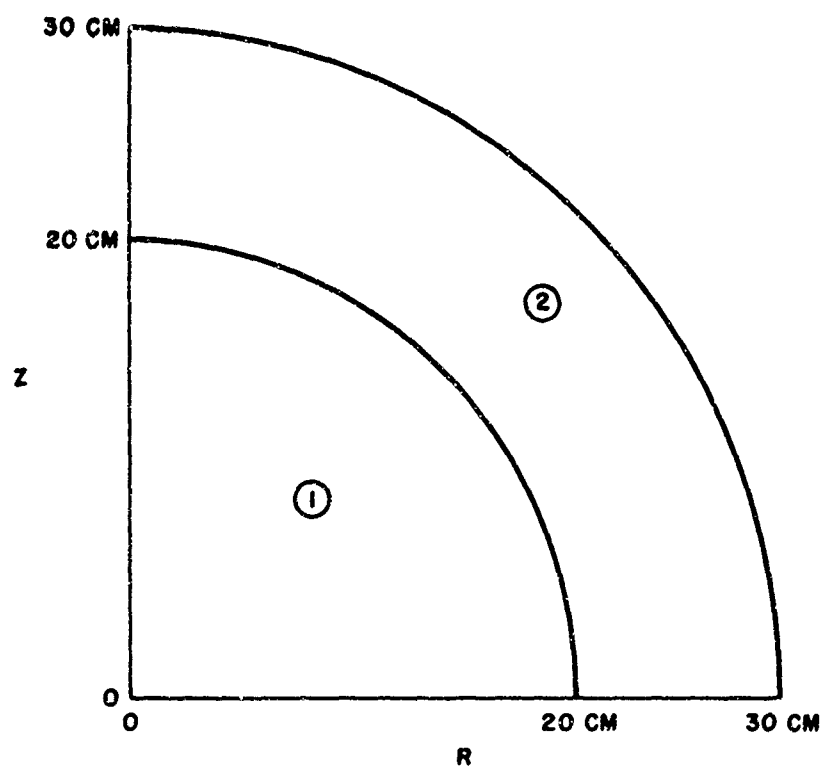


Fig. 15



- REGION ① $\rho = 0.8, \gamma = 5/3$
 $I = \text{SPECIFIC INTERNAL ENERGY}$
 $= 9 \times 10^{13} \text{ ERGS}$
 $U = V = 0$
- REGION ② $\rho = 1.2, \gamma = 5/3$
 $I = U = V = 0$

Fig. 16

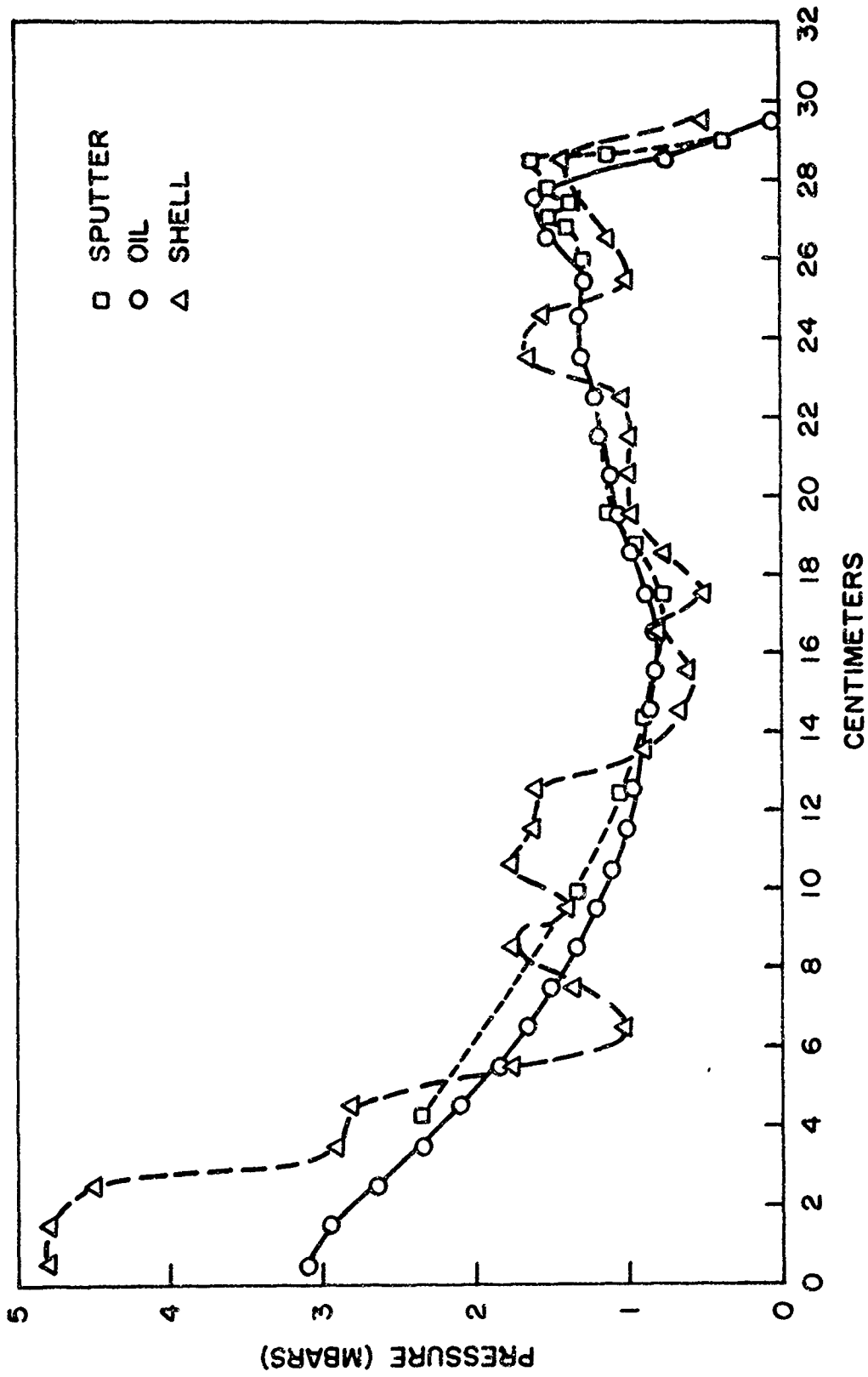


Fig. 17

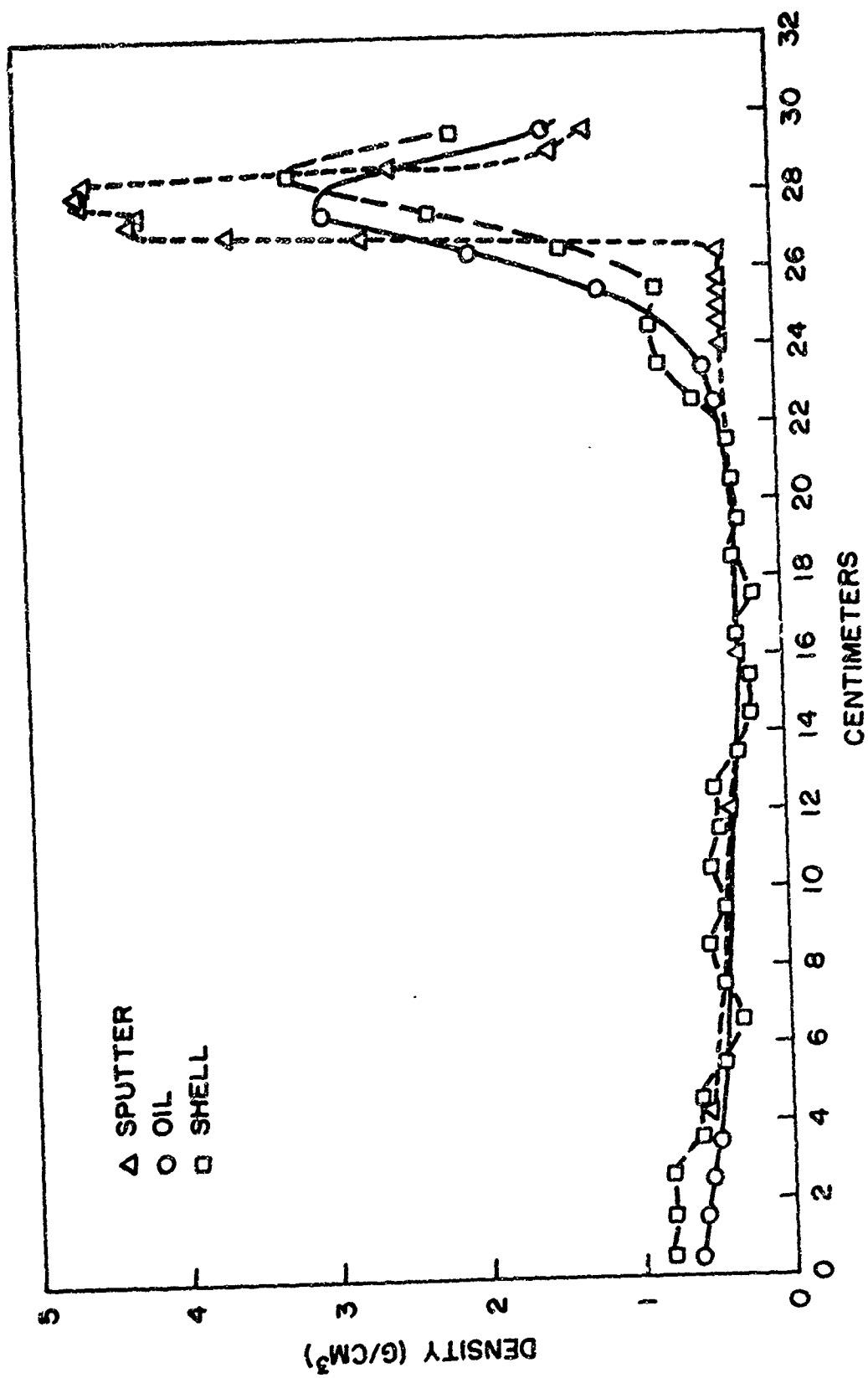


Fig. 18

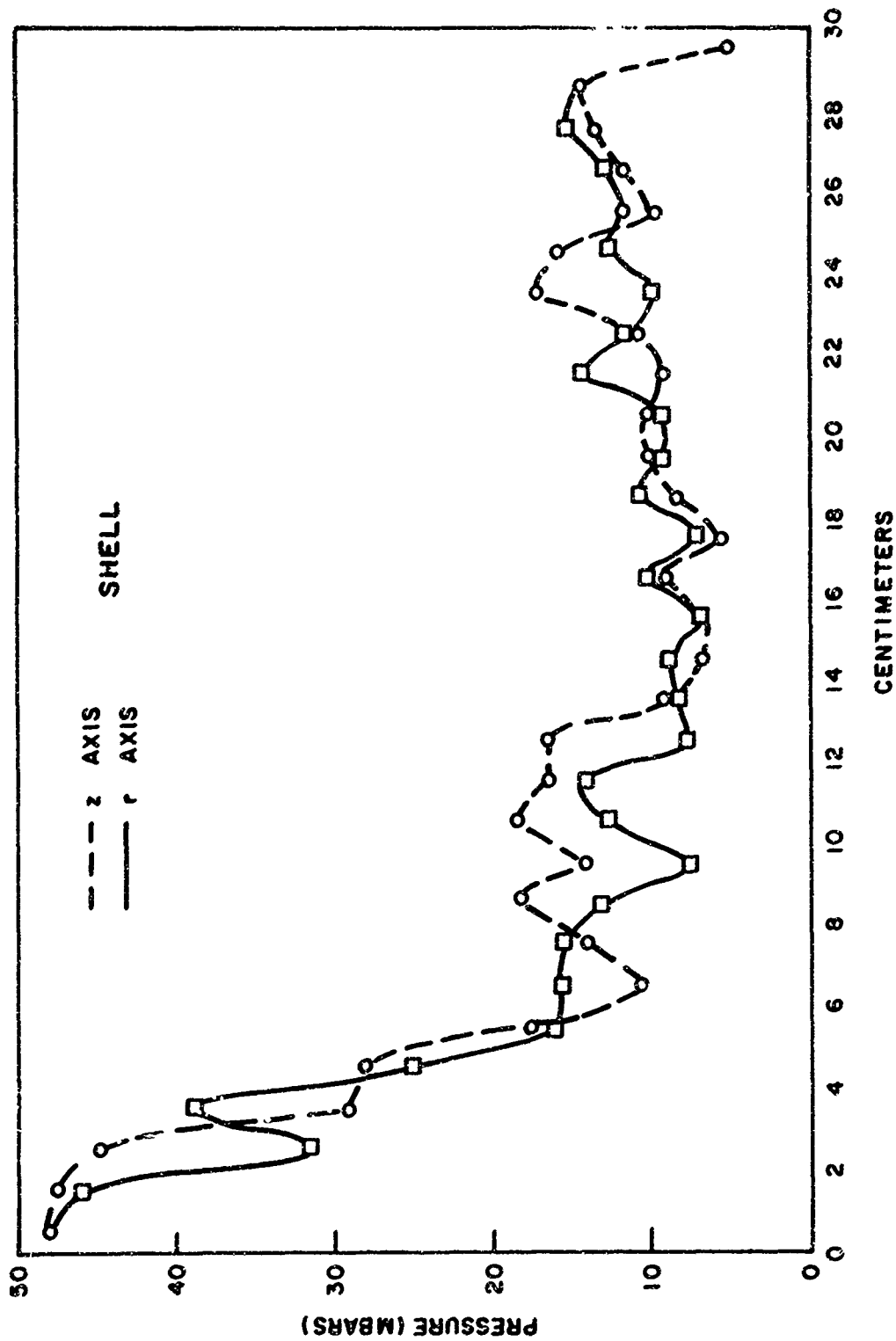


Fig. 19

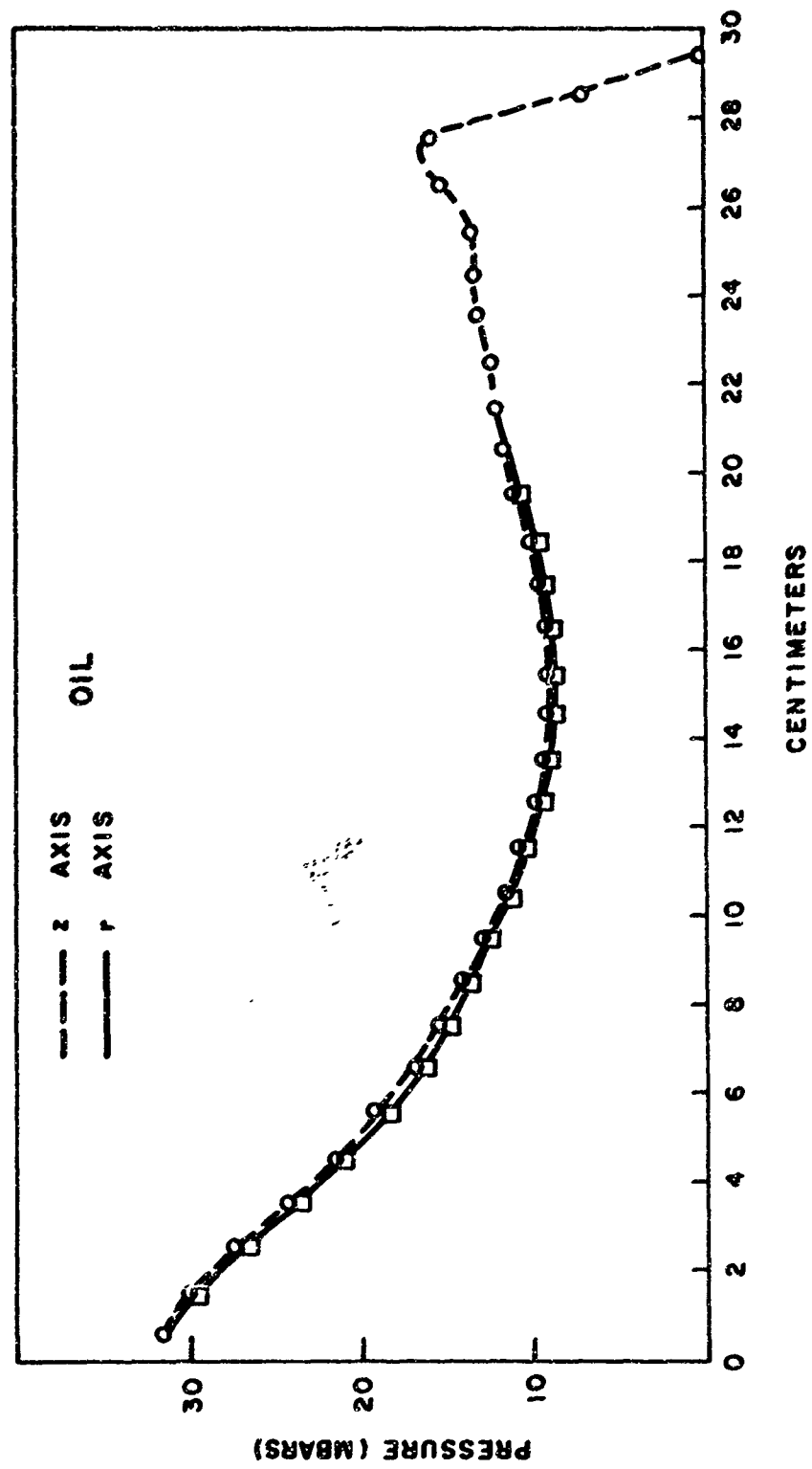


Fig. 20

5.1. FORTRAN IV LISTINGS OF CLAM

72.

C NOTE, THE BELOW SET OF
C DIMENSION, EQUIVALENCE AND COMMON IS
C TO BE USED WITH ALL SUBROUTINES IN
C CLAM, WITH THE EXCEPTION OF MAIN ROUTINE.
C D I M E N S I O N

DIMENSION AIX(4300),AM(130),
IAMX(4300),DX(52),ENDU(2),ITAB(502),IZ(100),
2RONE(2),TAB(502),TAB1(20),TAB1Y(21),TABR(20),
3TABUV(20),TABX(21),TABY(21),TAB152),TEMP(13),
4U(4300),V(4300),X(52),XL(130),XX(54),
5YL(130),Z(15),Y(100),YY(101),OY(100)

DIMENSION Iw1(130),Iw2(130)

COMMON	Z	,XX	,TAB	,YY		
COMMON	AID	,AIMAX	,AIX	,AJMAX	,AM	,AMD
COMMON	AMX	,DX	,ENDU	,FMX	,GXH	,GXX
COMMON	GYN	,GYX	,I	,IA	,IB	,IBA
COMMON	IBB	,IB	,IG	,II	,IIC	,IJ
COMMON	IR	,IRC	,IOV	,IOVL	,IWS	,IWSA
COMMON	IWSB	,IX	,IXN	,IXX	,IYN	,IYX
COMMON	J	,JA	,JI	,JIX	,K	,KE
COMMON	KF	,KK	,L	,LA	,LB	,LD
COMMON	LE	,LI	,LX	,M	,MI	,MIJ
COMMON	MJ	,AM	,MNP	,MX	,MXA	,MXS
COMMON	NZ	,NPKS	,NPP	,NT	,NX	,NY
COMMON	QOCCFL,RHO	,PONE,SLA,SLB		,TAB1		
COMMON	TAB1Y	,TABR	,TABUV	,TABX	,TABY	,TAM
COMMON	TAD	,TEMP	,TFMX	,TPIDY	,TX	,TY
COMMON	U	,V	,NPIDY	,WS	,WSA	,WSB
COMMON	WSC	,WSD	,WSE	,WSF	,WSG	,WSI
COMMON	WSL	,WSU	,WSV	,WSX	,WSY	,WS5
COMMON	XC	,XL	,YC	,YL	,YMAX	,WSR
COMMON	PE	,PM	,ITX	,TIY	,LF	,E
COMMON	PEE	,NPRK	,NYY	,DY	,NK	,SWITCH
COMMON	Iw1	,Iw2				

INPU0020
INPU0030
INPU0040
INPU0050
INPU0060
INPU0070
INPU0080
INPU0090
INPU0100
INPU0110
INPU0120
INPU0130
INPU0140
INPU0150
INPU0160
INPU0170
INPU0180
INPU0190
INPU0200
INPU0210
INPU0220
INPU0230
INPU0240
INPU0250
INPU0260
INPU0270
INPU0280
INPU0290
INPU0300
INPU0310
INPU0320
INPU0400
INPU0410
INPU0420

E Q U I V A L E N C E

OEQUIVALENCE	(Z,IZ,PROB),	(Z(2),CYCLE),	(Z(3),DT),
1(Z(4),PRINTS),	(Z(5),PRINTL),	(Z(6),DUMPT7),	(Z(7),CSTUP),
2(Z(8),PIDY),	(Z(9),TMZ),	(Z(10),GAM),	(Z(11),GAMU),
3(Z(12),GAMX),	(Z(13),ETH),	(Z(14),FFA),	(Z(15),FFB),
4(Z(16),TMDZ),	(Z(17),TMXZ),	(Z(18),XMAX),	(Z(19),TXMAX),
5(Z(20),TYMAX),	(Z(21),AMDM),	(Z(22),AMXM),	(Z(23),DNN),
6(Z(24),DMIN),	(Z(25),FEF),	(Z(26),DTNA),	(Z(27),CVIS),
7(Z(28),NPK),	(Z(29),NPRT),	(Z(30),NC),	(Z(31),NPC),
8(Z(32),NRC),	(Z(33),IMAX),	(Z(34),IMAXA),	(Z(35),JMAX),
9(Z(36),JMAXA),	(Z(37),KMAX),	(Z(38),KMAXA),	(Z(39),NMAX),
OEQUIVALENCE	(Z(40),ND),	(Z(41),KDT),	(Z(42),IXMAX),
1(Z(43),NOD),	(Z(44),NUPR),	(Z(45),NIMAX),	(Z(46),NJMAX),

INPU0430
INPU0440
INPU0450
INPU0460
INPU0470
INPU0480
INPU0490
INPU0500
INPU0510
INPU0520
INPU0530
INPU0540

2(Z(47),I1),	(Z(48),I2),	(Z(49),I3),	(Z(50),I4),	INPU0550
3(Z(51),N1),	(Z(52),N2),	(Z(53),N3),	(Z(54),N4),	INPU0560
4(Z(55),N5),	(Z(56),N6),	(Z(57),N7),	(Z(58),N8),	INPU0570
5(Z(59),N9),	(Z(60),N10),	(Z(61),N11),	(Z(62),NRM),	INPU0580
6(Z(63),TRAD),	(Z(64),XNRG),	(Z(65),SN),	(Z(66),DXN),	INPU0590
7(Z(67),RADER),	(Z(68),RADET),	(Z(69),RADEB),	(Z(70),DTRAD),	INPU0600
8(Z(71),REZFCT),	(Z(72),RSTOP),	(Z(73),SHELL),	(Z(74),BBOUND),	INPU0610
9(Z(75),TOZONE),	(Z(76),ECK),	(Z(77),SBOUND),	(Z(78),X1)	INPU0620
OEQUIVALENCE	(Z(79),X2),	(Z(80),Y1),	(Z(81),Y2),	INPU0630
1(Z(82),CABLH),	(Z(83),VISC),	(Z(84),T),	(Z(85),GMAX),	INPU0640
2(Z(86),MSGD),	(Z(87),MSGX),	(Z(88),GMADR),	(Z(89),GMAXR),	INPU0650
3(Z(90),S1),	(Z(91),S2),	(Z(92),S3),	(Z(93),S4),	INPU0660
4(Z(94),S5),	(Z(95),S6),	(Z(96),S7),	(Z(97),S8),	INPU0670
5(Z(98),S9),	(Z(99),S10)			INPU0680
OEQUIVALENCE	(Z,I2),	(XX(2),X(1)),	(TAB,ITAB),	INPU0690
1(Y(2),Y(1))				INPU0700
				INPU0710

C

NOTE . ALTHOUGH THE DIMENSIONS FOR THE CELL
QUANTITIES ARE 4300 IN CLAH, THE DIMENSIONS
FOR OIL ARE 3500, THAT IS, KEEP (IMAX)(JHAX)
+1 LESS THAN 3499.

74.

\$IBFTC MAIN LIST,DECK,REF

CMAIN

C L A M

***** M A I N *****

MAIN0010

MAIN0020

MAIN0030

C
C ** IF PROBLEM NUMBER IS NEGATIVE,
C CLAM WILL WRITE THE PARTICLES ON TAPE,
C PREPARING IT FOR A PIC RUN.
C *** NOTE (1 MATERIAL ONLY (1X))

MAIN0040

MAIN0050

MAIN0060

C CALL SLITE (0)
C INPUT ROUTINE CALCULATES THE ACTUAL GRID,
C DIMENSIONS AND INDICES.

10 CALL INPUT

MAIN0070

C PH1, READS IN DATA CARDS FOR THE
C PACKAGES, PH2 CALCULATES THE GEOMETRICS,
C PH3 THE PARTICLES, PH4 CALLS THE
C 6 POSSIBLE FITS THAT CALCULATE THE
C DENSITY, VELOCITIES AND INTERNAL ENERGY
C OF THE PARTICLES.

20 CALL PH1

MAIN0080

C OUTPUT CALCULATES THE VELOCITY (BOTH
C RADIAL AND AXIAL) AND SPECIFIC INTERNAL
C ENERGY OF EACH CELL FROM THE
C TOTAL MOMENTA AND INTERNAL
C ENERGY AND MASS OF EACH CELL.
C OUTPUT ALSO PREPARES A DUMP TAPE
C WHICH IS USED THEN TO START OIL.

30 CALL OUTPUT

MAIN0090

CALL EXIT

MAIN0100

END

MAIN0110

\$IBFIC INPUT LIST,DECK,REF	
SUBROUTINE INPUT	
C	INPU0910
C	INPU0730
C	INPU0940
C ***** NOTE (1 MATERIAL ONLY ((X))	INPU0950
MZ=150	INPU0960
C CLEAR Z BLOCK.	
DO 30 I=1,MZ	INPU0970
30 Z(I)=0.0	INPU0980
C READ IN HEADING CARD	INPU0990
READ (5,8012)IMS	INPU1000
IMS=1	INPU1010
WRITE (6,8012)(IMS)	INPU1020
WRITE (6,8100)	INPU1030
C READ IN PROBLEM CONSTANTS	INPU1040
C PROB=PROBLEM NO. AIMAX=IMAX,	
C AJMAX=JMAX, QCOOFL IS NOT USED-SET	
C TO ZERO, SHELL SET=2.,S8,S9 ARE	
C ZERO, SET N7 TO=TAPE NO.	
READ (5,8004)PROB,AIMAX,AJMAX,QCOOFL,SHELL,S8,S9,N7	INPU1050
IF(N7)40,40,50	INPU1060
40 N7=7	
50 CONTINUE	INPU1080
C MAX. NUMBER OF ZONES IN R DIRECTION.	
MI=50	INPU1090
C MAX. NUMBER OF ZONES IN Z DIRECTION.	
MJ=100	INPU1100
C MAX. NUMBER OF PARTICLES/CELL.	
MNP=400	INPU1110
C SIZE OF TABLE (TAB)	INPU1120
JTM=500	INPU1130
C MAXIMUM I*J	INPU1140
C MAX. NUMBER OF CELLS.	
60 MIJ=4299	INPU1150
C CALCULATE ADDITIONAL INDICES FOR CLAM AND OIL.	
70 IMAX=AIMAX	INPU1160
JMAX=AJMAX	INPU1170
IMAXA=IMAX+1	INPU1180
IXMAX=IMAXA+1	INPU1190
JMAXA=JMAX+1	INPU1200
KMAX=(IMAX*JMAX)+1	INPU1210
KMAXA=KMAX+1	INPU1220
WRITE (6,6048)(PROB,IMAX,JMAX)	INPU1230
C CHECK INPUT NOS. CONCERNED WITH GRID SIZE.	
101 IF(IMAX-MI)102,102,9901	INPU1240
102 IF(JMAX-MJ)104,104,9902	INPU1250
104 IF(KMAX-MIJ-1)106,106,9903	INPU1260
106 NOD=1	INPU1270
NPC=1	INPU1280
NRC=0	INPU1290

C	READ IN DX AND DY	INPU1300
	I=0	INPU1310
	J=0	IN 1320
	X(I)=0.0	I 1330
	Y(J)=0.0	INPU1340
2000	READ (5,8102) IWSA, IWSB, N1, N2, N3, N4, (TEMP(K), K=1, 4)	INPU1350
	L=1	INPU1360
C	COUNT NO. OF DIFFERENT DX OR DY.	INPU1370
	IF(N4) 2003, 2001, 2003	INPU1380
2001	IF(N3) 2004, 2002, 2004	INPU1390
2002	IF(N2) 2006, 2008, 2006	INPU1400
2003	L=L+1	INPU1410
2004	L=L+1	INPU1420
2006	L=L+1	INPU1430
2008	IF(IWSB) 2010, 2010, 2030	
C	PROCESS THE DX AND DY VALUES.	INPU1440
2010	DO 2014 N=1, L	INPU1450
	NK=IZ(N+50)	INPU1460
	DO 2012 K=1, NK	INPU1470
	I=I+1	INPU1480
	DX(I)=TEMP(N)	INPU1490
	X(I)=X(I-1)+DX(I)	INPU1500
2012	CONTINUE	INPU1510
2014	CONTINUE	INPU1520
	GO TO 2050	INPU1530
C	CALC THE Y AND DY VALUES	INPU1540
2030	DO 2034 N=1, L	INPU1550
	NK=IZ(N+50)	INPU1560
	DO 2032 K=1, NK	INPU1570
	J=J+1	INPU1580
	DY(J)=TEMP(N)	INPU1590
	Y(J)=Y(J-1)+DY(J)	INPU1600
2032	CONTINUE	INPU1610
2034	CONTINUE	INPU1620
2050	IF(IWSA) 2052, 2000, 2052	
C	IF(=) READ MORE DX OR DY DATA CARDS.	INPU1630
2052	IF(J-JMAX) 9905, 2053, 9905	
C	CHECK INPUT NUMBERS.	INPU1640
2053	IF(I-IMAX) 9906, 2054, 9906	INPU1650
2054	CONTINUE	INPU1660
	READ (5,8004) WS, WSA, WSB, SWITCH	
C	N1, AND N2 ARE THE 2 SCRATCH TAPES.	INPU1670
	N1=WS	INPU1680
	N2=WSA	INPU1690
	REWIND N1	INPU1700
	REWIND N2	
C	N4=MAX. NUMBER OF PARTICLES-1 PER RECORD.	INPU1710
	N4=WSB	INPU1720
	NPRI=N4	INPU1730
	NPRR=N4	

WRITE (6,6064)IMAX,(X(I),I=0,IMAX)	INPU1740
WRITE (6,6065)JMAX,(Y(J),J=0,JMAX)	INPU1750
MS=3.1415927	INPU1760
WSA=0.0	INPU1770
C CALCULATE THE AREA'S (TAU)=PI(R(I)**2-	
C R(I-1)**2).	
DO 1008 I=1,IMAX	INPU1780
WSB=WSA	INPU1790
WSA=X(I)**2	INPU1800
1008 TAU(I)=MS*(WSA-WSB)	INPU1810
C WRITE OUT X,Y,DX,DY, AND TAU VALUES.	
WRITE (6,6066)IMAX,(DX(I),I=1,IMAX)	INPU1820
WRITE (6,6067)JMAX,(DY(J),J=1,JMAX)	INPU1830
WRITE (6,6092)(IMAX,(TAU(I),I=1,IMAX))	INPU1840
1010 XMAX=X(IMAX)	INPU1850
TXMAX=XMAX*2.0	INPU1860
YMAX=Y(JMAX)	INPU1870
TYMAX=YMAX*2.0	INPU1880
C PIDY IS REALLY PI(3.1415927).	
PIDY=MS	INPU1890
C SET VELOCITIES, INTERNAL ENERGIES AND MASSES	
C TO 0.	
DO 1014 I=1,KMAX	INPU1910
U(I)=0.0	INPU1920
V(I)=0.0	INPU1930
AIX(I)=0.0	INPU1940
AMX(I)=0.0	INPU1950
1014 CONTINUE	INPU1960
C SET TOTAL ENERGY TO ZERO.	
ETH=0.0	INPU1970
C INITIALIZE MIN. MASS PARTICLE TO A LARGE NO.	
AMDM=1.E+28	INPU1980
AMXM=AMDM	INPU1990
GO TO 2016	INPU2000
C ERROR	INPU2010
9901 NK=101	INPU2020
GO TO 9999	INPU2030
9902 NK=102	INPU2040
GO TO 9999	INPU2050
9903 NK=104	INPU2060
GO TO 9999	INPU2070
C JMAX DOES NOT EQUAL THE SUM OF THE INPUT J	INPU2080
9905 NK=2052	INPU2090
GO TO 9999	INPU2100
C IMAX DOES NOT EQUAL THE SUM OF THE INPUT I	INPU2110
9906 NK=2053	INPU2120
9999 WRITE (6,8888)NK,I,J,X,L,M,N	INPU2130
PRINT 8888,NK,I,J,X,L,M,N	INPU2140
CALL DUMP	INPU2150
2016 RETURN	INPU2160

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C	FORMATS	
8004	FORMAT(7E10.5,I2)	INPU2170
80120	FORMAT (I1,7I1H) THIS IS THE CLAM PROGRAM AND THERE IS AN ERROR.	INPU2180
1)	INPU2190
8048	FORMAT(1H /9H PROB NO.F9.3,I2X,2HI=I2,2EX,2HJ=I2)	INPU2200
8064	FORMAT(1H /10H X(I) I=0,I2/(5F16.6))	INPU2210
8065	FORMAT(1H /10H Y(J) J=0,I2/(5F16.6))	INPU2220
8066	FORMAT(1H /11H DX(I) I=1,I2/(5F16.6))	INPU2230
8067	FORMAT(1H /11H DY(J) J=1,I2/(5F16.6))	INPU2240
8092	FORMAT(1H /13H AREA(I) I=1,I2/(5F16.6))	INPU2250
8100	FORMAT(1H /14H (SHELL INPUT))	INPU2260
8102	FORMAT(2I1,4I2,4E10.4)	INPU2270
8888	FORMAT(1H+/26H1 INPUT ERROR IN STATEMENT I5,I2X,I2H INDICES ARE 6 I 7)	INPU2280
END		INPU2290
		INPU2300

\$IBFTC PH1 LIST,DECK,REF
SUBROUTINE PH1

C		PH1 0010
C	***** NOTE (1 MATERIAL ONLY ((X))	PH1 0740
C	READ IN GEOMETRY ETC.	PH1 0950
	NPP=7	PH1 0960
	NPR=NPP-1	PH1 0980
	TPIDY=PIDY*2.0	PH1 0990
	ND=0	PH1 1000
	NX=0	PH1 1010
	NT=1	PH1 1020
	NY=1	PH1 1030
C	FIRST CARD OF EACH PACKAGE.	PH1 1040
	READ (5,8008)IX,LX,MX,TEMP(1),TEMP(2),TEMP(3)	
C	INITIALIZE THE NUMBER OF PACKAGES TO 0.	PH1 1050
	NPKG=0	
2015	IF(IX-1)9901,2018,2018	PH1 1060
2016	IX=I	PH1 1070
	LX=L	PH1 1080
	MX=M	PH1 1090
C	IF THERE ARE NO MORE PACKAGES GO COMPUTE TOTAL VALUES	PH1 1100
C	THE LAST CARD HAS A 2 PUNCH IN COL 1.	PH1 1110
2017	IF(IX-2)2018,7000,9902	PH1 1120
2018	J=0	PH1 1130
	NPKG=NPKG+1	PH1 1140
C	SET PACKAGE MASS AND ENERGY TO 0.	
	PE=0.0	PH1 1150
	PM=0.0	PH1 1160
C	ORIGIN FOR THE RADIUS VECTORS TO BE USED	
C	FOR THE FIT ROUTINES(1 THRU 6).	
	VC=TEMP(1)	PH1 1170
	XC=TEMP(2)	PH1 1180
C	S8 CONTAINS THE FIT NUMBER FOR THE	
C	PACKAGE IN QUESTION.	
	S8=TEMP(3)	PH1 1190
	WRITE (6,8100)(NPKG,MX)	PH1 1200
C	NOW READ IN THE GEOMETRY AND DENSITY,	
C	ENERGY AND VELOCITY CARDS.	
2020	READ (5,8008)I,L,M,(TEMP(N),N=1,6)	PH1 1210
	IWS=1	PH1 1220
	IF(I-5)2021,2040,2022	PH1 1230
C	IF=, THIS IS A RHO, VELOCITY OR ENERGY CARD.	
C	IF LESS, YOU HAVE READ ALL CARDS FOR THIS	
C	PACKAGE IN, PLUS THE FIRST CARD FROM THE	
C	NEXT PACKAGE.	
2021	IF(I-3)2060,9903,2026	PH1 1240
C	IF GREATER, EITHER A TRIANGLE OR PERTURBED ELLIPSE.	
2022	IF(L)9904,2030,2024	PH1 1250
C	A PERTURBED ELLIPSE.	
2024	IWS=7	PH1 1260

30.

GO TO 2030	PH1 1270
2026 IWS=3	PH1 1280
2027 IF(L)9905,2030,2028	PH1 1290
2028 IWS=5	PH1 1300
C A TRIANGLE.	
2030 IF(M)9906,2034,2032	PH1 1310
C IF=, DELETE THIS GEOMETRY.	
2032 IWS=IWS+1	PH1 1320
2034 J=J+1	PH1 1330
C TAB STORAGE CONTAINS THE COORDINATES OF	
C GEOMETRY.	
ITAB(J)=IWS	PH1 1340
DO 2036 N=1,NPR	PH1 1350
J=J+1	PH1 1360
2036 TAB(J)=TEMP(N)	PH1 1370
GO TO 2020	PH1 1380
C ONE ONLY RHO,I,U OR V ALLOWED PER PACKAGE	PH1 1390
C IF=, THIS IS A DENSITY CARD.	
2040 IF(L-1)9907,2046,2042	PH1 1400
C IF GREATER, EITHER A VELOCITY OR ENERGY CARD.	
2042 IF(L-3)2052,2058,9908	PH1 1410
C IF=, THIS IS A VELOCITY CARD, IF LESS, THIS IS A	
C ENERGY CARD.	
C DENSITY	PH1 1420
2046 DO 2048 N=1,6	PH1 1430
2048 TABR(N)=TEMP(N)	PH1 1440
GO TO 2020	PH1 1450
C ENERGY	PH1 1460
2052 DO 2054 N=1,6	PH1 1470
2054 TABI(N)=TEMP(N)	PH1 1480
GO TO 2020	PH1 1490
C VELOCITY (U AND V)	PH1 1500
2058 DO 2059 N=1,6	PH1 1510
2059 TABUV(N)=TEMP(N)	PH1 1520
GO TO 2020	PH1 1530
C OUTPUT DENSITY, ENERGY, AND VELOCITY PARAMETERS	PH1 1540
C ALL CARDS FOR THIS PACKAGE HAVE	
C BEEN READ IN.	
2060 IF(J-JTM)2070,2070,9915	PH1 1550
NO. OF PACKAGES EXCEED (72), NOTE	
C JTM SET=TO 500 IN INPUT, THUS MAX.	
C NO. OF PACKAGES = 72, UNLESS DIMENSIONS	
C ARE CHANGED.	
2070 WRITE (6,8036)(TABR(II),II=1,6)	PH1 1560
WRITE (6,8038)(TABI(II),II=1,6)	PH1 1570
WRITE (6,8040)(TABUV(II),II=1,6)	PH1 1580
C COMPUTE BOUNDARIES OF GEOMETRIES FOR EFFICIENCY IN	PH1 1590
C GENERATING OR DELETING PARTICLES	PH1 1600
2000 CALL PH2	PH1 1610
C COMPUTE I(0),I(N),J(0) AND J(N),FROM PREVIOUSLY	PH1 1620

C	COMPUTED VALUES, FOR UPPER AND LOWER LIMITS IN	PH1 1630
C	THE CELL MESH SCAN	PH1 1640
C	IXN=MINIMUM (I) OF GEOMETRY OF PACKAGE	
C	IYN=MINIMUM (J) OF GEOMETRY OF PACKAGE	
C	IXX=MAXIMUM (I) OF GEOMETRY OF PACKAGE	
C	IYX=MAXIMUM (J) OF GEOMETRY OF PACKAGE	
3001	IXN=1	PH1 1650
	IXX=1	PH1 1660
	IWS=IMAX-1	PH1 1670
3800	IF(IWS)9929,3820,3801	PH1 1680
3801	DO 3808 N=1,IWS	PH1 1690
	IF(X(N)-GXN)3802,3804,3804	PH1 1700
3802	IXN=IXN+1	PH1 1710
3804	IF(X(N)-GXX)3806,3806,3808	PH1 1720
3806	IXX=IXX+1	PH1 1730
3808	CONTINUE	PH1 1740
	IF(IXN)3812,3812,3814	PH1 1750
3812	IXN=1	PH1 1760
3814	IF(IMAX-IXX)3816,3818,3818	PH1 1770
3816	IXX=IMAX	PH1 1780
3818	IF(IXN-IXX)3820,3820,9930	PH1 1790
3820	IYN=1	PH1 1800
	IYX=1	PH1 1810
	IWS=JMAX-1	PH1 1820
3821	IF(IWS)9929,3834,3822	PH1 1830
3822	DO 3813 N=1,IWS	PH1 1840
3823	IF(Y(N)-GYN)3819,3817,3817	PH1 1850
3819	IYN=IYN+1	PH1 1860
3817	IF(Y(N)-GYX)3815,3815,3813	PH1 1870
3815	IYX=IYX+1	PH1 1880
3813	CONTINUE	PH1 1890
	IF(IYN)3824,3824,3826	PH1 1900
3824	IYN=1	PH1 1910
3826	IF(JMAX-IYX)3828,3830,3830	PH1 1920
3828	IYX=JMAX	PH1 1930
3830	IF(IYN-IYX)3834,3834,9931	PH1 1940
3834	WRITE (6,8044)IXN,IYN,IXX,IYX	PH1 1950
C	SCAN CELL MESH TO DETERMINE IF PARTICLES ARE TO BE	PH1 1960
C	GENERATED OR DELETED	PH1 1970
C	GENERATE PARTICLES	PH1 1980
4000	CALL PH3	PH1 1990
C	REARRANGE X,Y AND M FOR PARTICLES IF NECESSARY	PH1 2000
6011	LA=NY-NT	PH1 2010
	IF(LX)9947,6020,6022	PH1 2020
6020	ND=ND+LA	PH1 2030
	GO TO 6024	PH1 2040
6022	NX=NX+LA	PH1 2050
6024	NT=NY	PH1 2060
	ETH=ETH+PE	PH1 2070
C	REPLACE 606760000000 BY-18115198976	PH1 2080

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WS=(-ABS(-18115198976))	PH1 2090
6026 IF(LX)9933,6028,6030	PH1 2100
C REPLACE 244663000000 BY 22125740032	PH1 2110
6028 WS= ABS(22125740032)	PH1 2120
6030 WRITE (6,8501)LA,WS,PE,PM	PH1 2130
C GO READ IN NEXT PACKAGE	PH1 2140
6050 GO TO 2016	PH1 2150
7000 NMAX=NT	PH1 2160
C NMAX=MAX. NUMBER OF PARTICLES+1.	
C YOU HAVE PROCESSED ALL PACKAGES, ALL	
C PARTICLES, NOW GO TO THE OUTPUT.	
IF(AM(2))4051,4050,4051	PH1 2170
4050 N3=NRC	PH1 2180
GO TO 4060	PH1 2190
4051 NRC=NRC+1	PH1 2200
N3=NRC	PH1 2210
C N3=NO. OF PARTICLE RECORDS OF	
C N4 WORDS.	
IF(PROB)4052,4052,4060	PH1 2220
4052 WRITE (N2)(AM(I),XL(I),YL(I),IW1(I),IW2(I),I=2,NPRI)	PH1 2230
4060 N6=NMAX-(N4-1)*(N3-1)	PH1 2240
NOPR=N3	PH1 2250
REWIND N2	PH1 2260
GO TO 10000	PH1 2270
C ERROR	PH1 2280
9901 NK=2015	PH1 2290
GO TO 9999	PH1 2300
9902 NK=2017	PH1 2310
GO TO 9999	PH1 2320
9903 NK=2021	PH1 2330
GO TO 9999	PH1 2340
9904 NK=2022	PH1 2350
GO TO 9999	PH1 2360
9905 NK=2027	PH1 2370
GO TO 9999	PH1 2380
9906 NK=2030	PH1 2390
GO TO 9999	PH1 2400
9907 NK=2040	PH1 2410
GO TO 9999	PH1 2420
9908 NK=2042	PH1 2430
GO TO 9999	PH1 2440
9915 NK=2060	PH1 2450
GO TO 9999	PH1 2460
9929 NK=3800	PH1 2470
GO TO 9999	PH1 2480
9930 NK=3818	PH1 2490
GO TO 9999	PH1 2500
9931 NK=3830	PH1 2510
GO TO 9999	PH1 2520
9933 NK=6026	PH1 2530

GO TO 9999	PH1 2540
9947 NK=6011	PH1 2550
9999 WRITE (6,8888)NK	PH1 2560
PRINT 8888,NK	PH1 2570
CALL DUMP	PH1 2580
10000 RETURN	PH1 2590
C	PH1 2600
FORMATS	PH1 2610
8008 FORMAT (2I1,I5,E13.5,5E10.5)	PH1 2620
8036 FORMAT(1H07X,8HDENSITY 9X,1P6E16.6)	PH1 2630
8038 FORMAT(1H07X,8HENERGY 9X,1P6E16.6)	PH1 2640
8040 FORMAT(1H07X,8HVELOCITY9X,1P6E16.6/1H0/)	PH1 2650
8044 FORMAT(1H /6H I(1)=I2,4X,5HJ(1)=I2,4X,5HI(N)=I2,4X,5HJ(N)=I2)	PH1 2660
81000FORMAT(1H0///12HOPACKAGE NO.13,I20,15H PARTICLES/CELL//33X,2HA114X	PH1 2670
1,2HA214X,2HA314X,2HA414X,2HA514X,2HA6)	PH1 2680
85010FORMAT(1H0/I28,2H (A3,11H) PARTICLES22X,4HPE =1PE12.6,16X,4HPM =E1	PH1 2690
12.6)	PH1 2700
8888 FORMAT(23H1PH1 ERROR IN STATEMENTI5)	PH1 2710
END	

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\$IBFTC PH2	LIST,DECK,REF	
	SUBROUTINE PH2	PH2 0010
	CALCULATE THE PACKAGE GEOMETRIES	PH2 0020
C		PH2 0740
C		PH2 0950
C		PH2 0960
C	GENERATING OR DELETING PARTICLES	PH2 0970
C	J=VALUE OF LAST COORDINATE READ IN.	
	JT=J	PH2 0980
C	INITIALIZE OUTER BOUNDARIES.	
	GXN=XMAX	PH2 0990
	GYN=YMAX	PH2 1000
	GXX=0.0	PH2 1010
	GYX=0.0	PH2 1020
C	NPP=7(SET IN PH1).	
	DO 3700 J=1,JT,NPP	PH2 1030
C	IWS STORED IN ITAB ARRAY IN PH1.	
C	IF IWS=2(A TRIANGLE),IF=4(A RECTANGLE),	
C	IF=6,A ELLIPSE OR CIRCLE. IF IWS=8,A	
C	PERTURBED ELLIPSE. IF IWS IS LESS THAN	
C	THESE VALUES, THE DEFINITION STILL HOLDS, BUT	
C	NOW DELETE THIS GEOMETRY.	
	KK=(ITAB(J)-1)/2	PH2 1040
3007	IF(KK)9919,3010,3008	PH2 1050
3008	IF(KK-2)3100,3200,3009	PH2 1060
3009	IF(KK-4)3400,9920,9920	PH2 1070
C	TRIANGLE	PH2 1080
C	VERTICES CAN BE INPUTED IN ANY ORDER,	
C	X COORDINATE FIRST.	
C	SEARCH FOR THE LARGEST X(WSE) AND	
C	SMALLEST X(WSD).	
C	FIND MAXIMUM(WSE) AND MINIMUM(WSD) X COORDINATE	PH2 1090
3010	IF(TAB(J+1)-TAB(J+3))3011,3012,3013	PH2 1100
3011	WSE=TAB(J+3)	PH2 1110
	WSD=TAB(J+1)	PH2 1120
	GO TO 3014	PH2 1130
3012	TAB(J+1)=TAB(J+1)*1.0000001+1.0E-8	PH2 1140
3013	WSE=TAB(J+1)	PH2 1150
	WSD=TAB(J+3)	PH2 1160
3014	IF(TAB(J+5)-WSD)3020,3019,3016	PH2 1170
3016	IF(TAB(J+5)-WSE)3024,3017,3018	PH2 1180
3017	TAB(J+5)=TAB(J+5)*1.0000001+1.0E-8	PH2 1190
3018	WSE=TAB(J+5)	PH2 1200
	GO TO 3024	PH2 1210
3019	TAB(J+5)=TAB(J+5)*0.9999999-1.0E-8	PH2 1220
3020	WSD=TAB(J+5)	PH2 1230
C	ARRANGE VERTICES IN ASCENDING ORDER	PH2 1240
3024	IF(TAB(J+2)-TAB(J+4))3036,3034,3038	PH2 1250
3034	TAB(J+2)=TAB(J+2)*1.0000001+1.0E-8	PH2 1260
	GO TO 3038	PH2 1270

3036	WSA=TAB(J+1)	PH2 1280
	WSB=TAB(J+2)	PH2 1290
	TAB(J+1)=TAB(J+3)	PH2 1300
	TAB(J+2)=TAB(J+4)	PH2 1310
	TAB(J+3)=WSA	PH2 1320
	TAB(J+4)=WSB	PH2 1330
3038	IF(TAB(J+4)-TAB(J+6)) 3042,3040,3044	PH2 1340
3040	TAB(J+6)=TAB(J+6)*0.9999999-1.0E-8	PH2 1350
	GO TO 3044	PH2 1360
3042	WSA=TAB(J+3)	PH2 1370
	WSB=TAB(J+4)	PH2 1380
	TAB(J+3)=TAB(J+5)	PH2 1390
	TAB(J+4)=TAB(J+6)	PH2 1400
	TAB(J+5)=WSA	PH2 1410
	TAB(J+6)=WSB	PH2 1420
	GO TO 3024	PH2 1430
C	WSF=MINIMUM VALUE OF Y	
C	WSG=MAXIMUM VALUE OF Y	
3044	WSF=TAB(J+6)	PH2 1440
	WSG=TAB(J+2)	PH2 1450
C	COMPUTE SLOPES	PH2 1460
	SLA=(TAB(J+4)-TAB(J+2))/(TAB(J+3)-TAB(J+1))	PH2 1470
	SLB=(TAB(J+6)-TAB(J+2))/(TAB(J+5)-TAB(J+1))	PH2 1480
3053	IF(SLA-SLB) 3054,9921,3058	PH2 1490
3054	IF(SLA) 3056,9922,3064	PH2 1500
3056	IF(SLB) 3064,9923,3062	PH2 1510
3058	IF(SLA) 3062,9924,3056	PH2 1520
3062	WSA=TAB(J+3)	PH2 1530
	WSB=TAB(J+4)	PH2 1540
	WSC=SLA	PH2 1550
	TAB(J+3)=TAB(J+5)	PH2 1560
	TAB(J+4)=TAB(J+6)	PH2 1570
	SLA=SLB	PH2 1580
	TAB(J+5)=WSA	PH2 1590
	TAB(J+6)=WSB	PH2 1600
	SLB=WSC	PH2 1610
3064	IF(TAB(J+3)-TAB(J+5)) 3066,9925,3068	PH2 1620
3066	ITAB(J)=ITAB(J)+2	PH2 1630
	IWS=ITAB(J)-3	PH2 1640
	GO TO 3069	PH2 1650
3068	IWS=ITAB(J)-1	PH2 1660
3069	KE=J+1	PH2 1670
	KF=KE+5	PH2 1680
C	REPLACE 272545000000 BY 25058082816	PH2 1690
	WS= ABS(25058082816)	PH2 1700
	IF(IWS) 3072,3070,3072	PH2 1710
C	REPLACE 242543000000 BY 21836333056	PH2 1720
3070	WS= ABS(21836333056)	PH2 1730
3072	WRITE (6,8016)WS,(TAB(N),N=KE,KF)	PH2 1740
	WS=TAB(J+2)-SLB*TAB(J+1)	PH2 1750

TAB(J+1)=TAB(J+2)-SLA*TAB(J+1)	PH2 1760
TAB(J+6)=(TAB(J+6)-TAB(J+4))/(TAB(J+5)-TAB(J+3))	PH2 1770
TAB(J+5)=TAB(J+4)-TAB(J+6)*TAB(J+3)	PH2 1780
TAB(J+2)=SLA	PH2 1790
TAB(J+3)=WS	PH2 1800
TAB(J+4)=SLB	PH2 1810
GO TO 3600	PH2 1820
C RECTANGLE	PH2 1830
3100 ITAB(J)=ITAB(J)+2	PH2 1840
IWS=ITAB(J)-5	PH2 1850
C REPLACE 272545000000 BY 25058082816	PH2 1860
WS= ABS(25058082816)	PH2 1870
IF(IWS)3110,3105,3110	PH2 1880
C REPLACE 242543000000 BY 21836333056	PH2 1890
3105 WS= ABS(21836333056)	PH2 1900
3110 WRITE (6,8020)WS,TAB(J+1),TAB(J+2),TAB(J+3),TAB(J+4)	PH2 1910
WSD=TAB(J+1)	PH2 1920
WSE=TAB(J+2)	PH2 1930
WSF=TAB(J+3)	PH2 1940
WSG=TAB(J+4)	PH2 1950
GO TO 3600	PH2 1960
C ELLIPSE OR CIRCLE	PH2 1970
3200 IF(ABS(TAB(J+1)-TAB(J+2))-1.0E-8)3300,3300,3202	PH2 1980
3202 IF(TAB(J+2))9926,3300,3203	PH2 1990
C ELLIPSE WITH NO PERTURBATION	PH2 2000
3203 ITAB(J)=ITAB(J)+2	PH2 2010
IWS=ITAB(J)-7	PH2 2020
C REPLACE 272545000000 BY 25058082816	PH2 2030
WS= ABS(25058082816)	PH2 2040
IF(IWS)3210,3205,3210	PH2 2050
C REPLACE 242543000000 BY 21836333056	PH2 2060
3205 WS= ABS(21836333056)	PH2 2070
3210 WRITE (6,8024)WS,TAB(J+1),TAB(J+2),TAB(J+3),TAB(J+4)	PH2 2080
3215 WSD=TAB(J+3)-TAB(J+1)	PH2 2090
WSE=TAB(J+3)+TAB(J+1)	PH2 2100
WSF=TAB(J+4)-TAB(J+2)	PH2 2110
WSG=TAB(J+4)+TAB(J+2)	PH2 2120
TAB(J+1)=TAB(J+1)**2	PH2 2130
TAB(J+2)=TAB(J+2)**2	PH2 2140
GO TO 3600	PH2 2150
C CIRCLE	PH2 2160
3300 ITAB(J)=ITAB(J)+4	PH2 2170
IWS=ITAB(J)-9	PH2 2180
TAB(J+2)=TAB(J+1)	PH2 2190
C REPLACE 272545000000 BY 25058082816	PH2 2200
WS= ABS(25058082816)	PH2 2210
IF(IWS)3310,3305,3310	PH2 2220
C REPLACE 242543000000 BY 21836333056	PH2 2230
3305 WS= ABS(21836333056)	PH2 2240
3310 WRITE (6,8028)WS,TAB(J+1),TAB(J+3),TAB(J+4)	PH2 2250

GO TO 3215	PH2 2260
C ELLIPSE WITH PERTURBATION	PH2 2270
3400 ITAB(J)=ITAB(J)+4	PH2 2280
WS=1.0-(TAB(J+5)/TAB(J+1))**2	PH2 2290
IWSA=ITAB(J+7)	PH2 2300
OTAB(J+7)=(TAB(J+6)-TAB(J+4)-TAB(J+2)*SQRT(WS))/	PH2 2310
1 ((TAB(J+5)*(TAB(J+5)-TAB(J+1)))**2)	PH2 2320
IWS=ITAB(J)-11	PH2 2330
KE=J+1	PH2 2340
KF=KE+6	PH2 2350
C REPLACE 272545000000 BY 25058082816	PH2 2360
WSA= ABS(25058082816)	PH2 2370
IF(IWS)3410,3405,3410	PH2 2380
C REPLACE 242543000000 BY 21836333056	PH2 2390
3405 WSA= ABS(21836333056)	PH2 2400
3410 WRITE (6,8032)WSA,(TAB(N),N=KE,KF)	PH2 2410
3415 IF(WS)9927,9927,3420	PH2 2420
3420 IF(TAB(J+3))9928,3425,9928	PH2 2430
3425 TAB(J+3)=TAB(J+7)	PH2 2440
ITAB(J+7)=IWSA	PH2 2450
WSA=TAB(J+2)+TAB(J+2)/4.0	PH2 2460
WSD=0.0	PH2 2470
WSE=TAB(J+1)+TAB(J+1)/4.0	PH2 2480
WSF=TAB(J+4)-WSA	PH2 2490
WSG=TAB(J+4)+WSA	PH2 2500
C DETERMINE BOUNDARIES OF GEOMETRIES	PH2 2510
3600 IF(WSD-GXN)3602,3604,3604	PH2 2520
C MAXIMUM (X)	
3602 GXN=WSD	PH2 2530
3604 IF(WSE-GXX)3608,3608,3606	PH2 2540
C MINIMUM (X)	
3606 GXX=WSE	PH2 2550
3608 IF(WSF-GYN)3610,3612,3612	PH2 2560
C MAXIMUM (Y)	
3610 GYN=WSF	PH2 2570
3612 IF(WSG-GYX)3700,3700,3614	PH2 2580
C MINIMUM (Y)	
3614 GYX=WSG	PH2 2590
3700 CONTINUE	PH2 2600
J=JT	PH2 2610
GO TO 10000	PH2 2620
C E R R O R	PH2 2630
9919 NK=3007	PH2 2640
GO TO 9999	PH2 2650
9920 NK=3009	PH2 2660
GO TO 9999	PH2 2670
9921 NK=3053	PH2 2680
GO TO 9999	PH2 2690
9922 NK=3054	PH2 2700
GO TO 9999	PH2 2710

88.

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9923 NK=3056
      GO TO 9999
9924 NK=3058
      GO TO 9999
9925 NK=3064
      GO TO 9999
9926 NK=3202
      GO TO 9999
9927 NK=3415
      GO TO 9999
9928 NK=3420
9999 WRITE (6,8888)NK
      PRINT 8888,NK
      CALL DUMP
10000 RETURN
8016 FORMAT(15H0TRIANGLE ---- A3,7H -----1P0E10.6)
8020 FORMAT(15H0RECTANGLE --- A3,7H -----1P6E10.6)
8024 FORMAT(15H0ELLIPSE ---- A3,7H -----1P6E10.6)
8028 FORMAT(15H0CIRCLE ----- A3,7H -----1PE10.0,10X,4E10.6)
8032 FORMAT(15H0P ELLIPSE --- A3,7H -----1P6E10.6)
8888 FORMAT(23H1PH2 ERROR IN STATEMENT15)
      END
```

PH2 2720
PH2 2730
PH2 2740
PH2 2750
PH2 2760
PH2 2770
PH2 2780
PH2 2790
PH2 2800
PH2 2810
PH2 2820
PH2 2830
PH2 2840
PH2 2850
PH2 2860
PH2 2870
PH2 2880
PH2 2890
PH2 2900
PH2 2910
PH2 2920
PH2 2930

\$IBFTC PH3	LIST,DECK,REF	
	SUBROUTINE PH3	PH3 0010
C	GENERATE (OR DELETE) THE PARTICLES	PH3 0020
C		PH3 0740
C		PH3 0950
C		PH3 0960
C		PH3 0970
C	SCAN CELL MESH TO DETERMINE IF PARTICLES ARE TO BE	PH3 0980
C	GENERATED OR DELETED	PH3 0990
C	GENERATE PARTICLES	PH3 1000
C	SAVE CURRENT VALUES OF COUNTERS.	
4000	IA=I	PH3 1010
	JA=J	PH3 1020
	IJ=K	PH3 1030
	JT=L	PH3 1040
	IF(IX-1)9932,4010,9932	PH3 1050
4010	IF(MX-MNP)4012,4012,9935	PH3 1060
C	IF GREATER, YOU TRIED TO GENERATE MORE THAN	
C	400 PARTICLES / CELL.	
4012	WS=MX	PH3 1070
	FMX=SQRT(WS)	PH3 1080
	MXS=FMX+.5	PH3 1090
4011	IF(MXS*MXS-MX)9936,4013,9936	PH3 1100
C	IF(GREATER OR LESS) THE NO. OF PARTICLES / CELL	
C	THAT YOU REQUESTED WAS NOT N SQ. WHERE	
C	N IS FROM 1 TO 20.	
4013	MXA=1-MX	PH3 1110
	TFMX=.5/FMX	PH3 1120
	WPIDY=TPIDY/FMX	PH3 1130
4015	IF(MXA)4018,4018,9937	PH3 1140
C	IF GREATER, YOU HAVE FAILED TO SPECIFY THE	
C	NO. OF PARTICLES TO GENERATE.	
4018	NY=NT	PH3 1150
	DO 5700 I=IXN,IXX	PH3 1160
C	COMPUTE THE COORDINATE OF THE PARTICLE	PH3 1170
C	UNDER CONSIDERATION	PH3 1180
	WS5=DX(I)/FMX	PH3 1190
C	THE VOLUME OF THE SUBDIVIDED CELL =	
C	PI(2.*XL(N)DY/N*DY/N).	
	TABX(1)=X(1)-TFMX*DX(1)	PH3 1200
4019	IF(MXA)4020,4024,9938	PH3 1210
4020	DO 4022 K=2,MXS	PH3 1220
C	WE START AT THE RIGHT AND TOP OF CELL(K).	
C	SET UP ARRAY FOR X COORDINATES OF THE	
C	PARTICLES.	
4022	TABX(K)=TABX(K-1)-WS5	PH3 1230
C	J LOOP, LIMITS OF Y FOR THIS PACKAGE.	
4024	DO 5700 J=IYN,IVX	PH3 1240
	TAM=WPIDY*KS5*DY(J)	PH3 1250
C	TAM= 2PI/N*DX/N*DY	

E=0.0	PH3 1260
IIWS=0	PH3 1270
IWS=0	PH3 1280
IB=0	PH3 1290
WS=DY(J)/FMX	PH3 1300
TABY(1)=Y(J)-TFMX*DY(J)	PH3 1310
C MXS=N	
DO 4026 K=2,MXS	PH3 1320
C SET UP ARRAY FOR Y COORDINATES OF THE	
C PARTICLES.	
4026 TABY(K)=TABY(K-1)-WS	PH3 1330
C K USED FOR THE CELL QUANTITIES.	
K=(J-1)*IMAX+1	PH3 1340
4028 IBB=IB/MXS	PH3 1350
IB=IB+1	PH3 1360
IBA=MOD(IB,MXS)	PH3 1370
C TX=X COORDINATE OF PARTICLE IN QUESTION.	
TX=TABX(1BA+1)	PH3 1380
C TY=Y COORDINATE OF PARTICLE IN QUESTION.	
TY=TABY(1BB+1)	PH3 1390
C GENERATE () DELETE THE PARTICLE	PH3 1400
ID=0	PH3 1410
IG=0	PH3 1420
DO 4200 L=1,JA,NPP	PH3 1430
KK=ITAB(L)	PH3 1440
IF(KK-5)4062,4078,4078	PH3 1450
C TRIANGLE	PH3 1460
4062 WSX=(TY-TAB(L+1))/TAB(L+2)	PH3 1470
IF(WSX-TX)4064,4064,4200	PH3 1480
4064 WSX=(TY-TAB(L+3))/TAB(L+4)	PH3 1490
IF(WSX-TX)4200,4066,4066	PH3 1500
4066 WSY=TAB(L+6)*TX+TAB(L+5)	PH3 1510
IF(KK-2)4068,4068,4072	PH3 1520
4068 IF(WSY-TY)4200,4070,4070	PH3 1530
4070 GO TO (4074,4076,4074,4076),KK	PH3 1540
4072 IF(WSY-TY)4070,4070,4200	PH3 1550
4074 ID=1	PH3 1560
GO TO 4200	PH3 1570
4076 IG=1	PH3 1580
GO TO 4200	PH3 1590
4078 KK=KK-4	PH3 1600
4077 IF(KK-8)4079,4094,9939	PH3 1610
4079 GO TO (4080,4080,4090,4090,4092,4092,4094),KK	PH3 1620
C RECTANGLE	PH3 1630
4080 IF(TAB(L+1)-TX)4082,4082,4200	PH3 1640
4082 IF(TAB(L+2)-TX)4200,4084,4084	PH3 1650
4084 IF(TAB(L+3)-TY)4086,4086,4200	PH3 1660
4086 IF(TAB(L+4)-TY)4200,4088,4088	PH3 1670
4088 GO TO (4074,4076),KK	PH3 1680
C ELLIPSE WITH NO PERTURBATION	PH3 1690

4090	KK=KK-2	PH3 1700
	IF((TX-TAB(L+3))*2/TAB(L+1)+(TY-TAB(L+4))*2	PH3 1710
	1/TAB(L+2)-1.0)4088,4088,4200	PH3 1720
C	CIRCLE	PH3 1730
4092	KK=KK-4	PH3 1740
	OIF((TX-TAB(L+3))*2+(TY-TAB(L+4))*2-TAB(L+1))	PH3 1750
	1 4088,4088,4200	PH3 1760
C	ELLIPSE WITH PERTURBATION	PH3 1770
4094	KK=KK-6	PH3 1780
	OIF((TX/TAB(L+1))*2+(TY-TAB(L+4)-TAB(L+3)*(TX*	PH3 1790
	1 (TX-TAB(L+1))*2)**2/TAB(L+2)-1.0)4088,4088,4200	PH3 1800
4200	CONTINUE	PH3 1810
C	IF ID=1 DELETE	PH3 1820
4201	IF(ID)9940,4310,4800	PH3 1830
C	IF ID=0 AND IG=0 DELETE	PH3 1840
4310	IF(IG)9941,4800,4312	PH3 1850
C	GENERATE PARTICLE	PH3 1860
4312	NY=NY+1	PH3 1870
	IF(IIWS)23,22,23	PH3 1880
22	IIWS=1	PH3 1890
23	IWS=1	PH3 1900
	NYY=NYY+1	PH3 1910
	CALL PH4	PH3 1950
C	RETURN FROM PH4 WITH THE FOLLOWING DATA,	
C	WSR=PARTICLE DENSITY	
C	WSI=PARTICLE SPECIFIC INTERNAL ENERGY	
C	WSU=RADIAL VELOCITY COMPONENT OF PARTICLE	
C	WSV=AXIAL VELOCITY COMPONENT OF PARTICLE	
4332	N=NYY	PH3 1960
	IF(IIWS)4335,4335,24	PH3 1970
24	IIWS=-1	PH3 1980
4333	IF(AMX(K))9951,4335,4334	PH3 1990
C	CALCULATE PACKAGE ENERGY.	
4334	E=((U(K)**2+V(K)**2)/(AMX(K)))*.5+AIX(K)	PH3 2000
C	SET THE PARTICLE COORDINATES INTO THE	
C	PROPER ARRAYS.	
4335	XL(N)=TX	PH3 2010
	YL(N)=TY	PH3 2020
C	SET I AND J OF CELL K(LOCATION OF PARTICLE).	
	IW1(N)=I	PH3 2030
	IW2(N)=J	PH3 2040
C	CALCULATE PARTICLE MASS AS	
C	=2PI/N*DX/N*DY*XL(N)*RHO.	
	AM(N)=TAM*TX*WSR	PH3 2050
4341	IF(LX)9945,4342,4344	PH3 2060
4342	WS=AM(N)*WSI	PH3 2070
	IF(AM(N)-ANDM)16,15,15	PH3 2080
16	ANDM=AM(N)	PH3 2090
15	CONTINUE	PH3 2100
	PM=PM+AM(N)	PH3 2110

AM(N)=-AM(N)	PH3 2120
GO TO 4346	PH3 2130
4344 WS=AM(N)*WSI	PH3 2140
IF 'AM(N)-AMXM)18,17,17	PH3 2150
18 AMXM=AM(N)	PH3 2160
17 AIX(K)=AIX(K)+WS	PH3 2170
PM=PM+AM(N)	PH3 2180
C SUM UP MASS, BOTH COMPONENTS OF MOMENTA	
C AND TOTAL INTERNAL ENERGY IN CELL K.	
AMX(K)=AMX(K)+AM(N)	PH3 2190
4346 U(K)=U(K)+ABS(AM(N))*WSU	PH3 2200
V(K)=V(K)+ABS(AM(N))*WSV	PH3 2210
IF(NY-NPRR)4800,14,9945	PH3 2220
14 NRC=NRC+1	PH3 2230
NPRR=NPRR+NPRI-1	PH3 2240
IF(PROB)5000,5000,5001	PH3 2250
C WRITE PARTICLES ON TAPE IF THIS IS TO	
C BE A PIC RUN.	
5000 WRITE (N2)(AM(N),XL(N),YL(N),IW1(N),IW2(N),N=2,NPRI)	PH3 2260
5001 NYY=1	PH3 2270
3 DO 2 N=2,NPRI	PH3 2280
C SET PARTICLE ARRAYS TO ZERO.	
XL(N)=0.0	PH3 2290
YL(N)=0.0	PH3 2300
AM(N)=0.0	PH3 2310
IW1(N)=0	PH3 2320
IW2(N)=0	PH3 2330
2 CONTINUE	PH3 2340
4800 IF(MX-2B)9946,4880,4028	PH3 2350
C A L C U L A T E E N E R G Y F O R P K G	PH3 2360
4880 IF(IWS)4900,5700,4900	PH3 2370
4900 IF(AMX(K))9951,5700,4910	PH3 2380
4910 PEE=(U(K)**2+V(K)**2)/(AMX(K))*0.5+AIX(K)	PH3 2390
4930 IF(E)4950,4940,4940	PH3 2400
4940 PEE=PEE-E	PH3 2410
4950 PE=PE+PEE	PH3 2420
5700 CONTINUE	PH3 2430
I=IA	PH3 2440
J=JA	PH3 2450
K=IJ	PH3 2460
L=JT	PH3 2470
GO TO 10000	PH3 2480
C E R R O R	PH3 2490
9932 NK=4000	PH3 2500
GO TO 9999	PH3 2510
9935 NK=4010	PH3 2520
GO TO 9999	PH3 2530
9936 NK=4011	PH3 2540
GO TO 9999	PH3 2550
9937 NK=4015	PH3 2560

GO TO 9999	PH3 2570
9938 NK=4019	PH3 2580
GO TO 9999	PH3 2590
9939 NK=4077	PH3 2600
GO TO 9999	PH3 2610
9940 NK=4201	PH3 2620
GO TO 9999	PH3 2630
9941 NK=4310	PH3 2640
GO TO 9999	PH3 2650
9945 NK=4341	PH3 2660
GO TO 9999	PH3 2670
9946 NK=4800	PH3 2680
GO TO 9999	PH3 2690
9951 NK=4905	PH3 2700
9999 WRITE (6,8888)NK,I,J,K,L,M,N	PH3 2710
PRINT 8888,NK,I,J,K,L,M,N	PH3 2720
CALL DUMP	PH3 2730
10000 RETURN	PH3 2740
8888 FORMAT(1H+/26H1 P H 3 ERROR IN STATEMENTI5,12X,12H INDICES ARE6I7)	PH3 2750
END	PH3 2760

94.

\$IBFTC PH4 LIST,DECK,REF
SUBROUTINE PH4

C
C
C THE ACTUAL COORDINATES USED IN THE FIT
C SUBROUTINES IS TTX=TX-XC,TTY=TY-YC.
TTX=TX-XC
TTY=TY-YC
LL=S8
GO TO(1,2,3,4,5,6),LL
1 CALL FIT1
GO TO 7
2 CALL FIT2
GO TO 7
3 CALL FIT3
GO TO 7
4 CALL FIT4
GO TO 7
5 CALL FIT5
GO TO 7
6 CALL FIT6
7 RETURN
END

PH4 0010
PH4 0730
PH4 0940

PH4 0950
PH4 0960
PH4 0970
PH4 0980
PH4 0990
PH4 1000
PH4 1010
PH4 1020
PH4 1030
PH4 1040
PH4 1050
PH4 1060
PH4 1070
PH4 1080
PH4 1090
PH4 1100
PH4 1110

SUBROUTINE PH4 DETERMINES WHICH ONE OF THE SIX POSSIBLE FITS TO CALL FOR, TO ASSIGN A DENSITY ENERGY, AND VELOCITIES TO PARTICLE N. THE FIT NUMBER IS SPECIFIED IN COLUMNS 31*40 OF THE FIRST CARD OF EACH PACKAGE(SEE SECTION 2.1)

\$IBFTC FIT1 LIST,DECK,REF
SUBROUTINE FIT1

C
C
C WS=SQRT(TTX**2+TTY**2)
C DENSITY
C WSR=TABR(1)+TABR(2)*(TTY-TABR(3))
C ENERGY
C WSI=TABI(1)+TABI(2)*(TTY-TABI(3))
C VELOCITIES
C WS=TABUV(1)+TABUV(2)* TTY-TABUV(3))
C WSU=0.0
C WSV=WS
C RETURN
C END

FIT10010
FIT10730
FIT10940
FIT10950
FIT10960
FIT10970
FIT10980
FIT10990
FIT11000
FIT11010
FIT11020
FIT11030
FIT11040
FIT11050

\$IBFTC FIT2 LIST,DECK,REF
SUBROUTINE FIT2

C
C
C WS=SQRT(TTX**2+TTY**2)
C DENSITY
C WSR=((TTX-TABR(1))/TABR(2))**2+((TTY-TABR(3))/
1TABR(4))**2
C ENERGY
C WSI=TABI(1)+TABI(2)*TTX+TABI(3)*TTX**2
C 1+TABI(4)*TTY+TABI(5)*TTY**2
C VELOCITIES
C WSV=TABUV(1)+TABUV(2)*TTY
C WSU=TABUV(3)+TABUV(4)*TTY
C RETURN
C END

FIT20010
FIT20730
FIT20940
FIT20950
FIT20960
FIT20970
FIT20980
FIT20990
FIT21000
FIT21010
FIT21020
FIT21030
FIT21040
FIT21050
FIT21060

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\$IBFTC FIT3 LIST,DECK,REF
SUBROUTINE FIT3

FIT30010
FIT30730
FIT30940

C
C
C THIS FIT FOR SIN KZ/KZ *****
WS=SQRT(TTX**2+TTY**2)
C DENSITY
WSR=TABR(1)+TABR(2)*{TTY-TABR(3)}
WSA=TTY/TABI(2)
WSB=WSA*PIDY*2.
WSC=SIN(WSB)
WSI=WSC/WSA*TABI(1)
WS=TABUV(1)+TABUV(2)*{TTY-TABUV(3)}
WSU=0.
WSV=WS
WSI=WSI*TABI(3)
C TABI(3) US SCALE FACTOR FOR YIELD NORMALLY SET TO 1.
RETURN
END

FIT30950
FIT30960

\$IBFTC FIT4 LIST,DECK,REF
SUBROUTINE FIT4
RETURN
END

FIT40010
FIT40020
FIT40030

\$IBFTC FIT5 LIST,DECK,REF
SUBROUTINE FIT5
RETURN
END

FIT50010
FIT50020
FIT50030

\$IBFTC FIT6 LIST,DECK,REF
SUBROUTINE FIT6
RETURN
END

FIT60010
FIT60020
FIT60030

\$IBFTC OUTPUT LIST,DECK,REF
SUBROUTINE OUTPUT

C C O M M O N

C L A M ***** O U T P U T *****

C

C

C NOTE (1 MATERIAL ONLY ((X))

C

C

C

PACKAGES HAVE BEEN READ IN AND PROCESSED
COMPUTE TOTAL ENERGIES AND TOTAL MASSES

E=ETH

WRITE (6,8104)

7001 ND=ND+1

IF(E)6000,6000,6001

6000 AMDM=0.0

AMXM=0.0

GO TO 7016

6001 AMDM=AMDM/2.0

AMXM=AMXM/2.0

7013 IF(AMDM)9901,9901,7014

7014 IF(AMXM)9902,9902,7016

7016 ETH=0.0

TMDZ=0.0

TMXZ=0.0

DO 7012 I=2,KMAX

7005 IF(AMX(I))9904,7012,7006

7006 CONTINUE

C SUM UP TOTAL (X) MASS IN GRID.

TMXZ=AMX(I)+TMXZ

C CALCULATE SPECIFIC INTERNAL ENERGY/CELL (K).

AIX(I)=AIX(I)/AMX(I)

7008 WS=AMX(I)

C CALCULATE RADIAL AND AXIAL VELOCITIES BY

C CONSERVING BOTH COMPONENTS OF MOMENTA.

U(I)=U(I)/WS

V(I)=V(I)/WS

C SUM UP TOTAL ENERGY IN SYSTEM.

ETH=ETH+((U(I)**2+V(I)**2)/2.0+AIX(I))*WS

GO TO 7012

7012 CONTINUE

TMZ=TMDZ+TMXZ

WRITE (6,8072)ETH,E,TMDZ,TMXZ,TMZ

IWS=ND-1

IWSA=NMAX-ND

IWSB=NMAX-1

WRITE (6,8073)(IWS,IWSA,IWSB)

C PUT INPUT ON BINARY TAPE 7

7113 REWIND N7

C

C

WRITE TAPE FOR OIL CODE.

OUTP0010

INPU0720

OUTP0020

OUTP0030

OUTP0750

OUTP0960

OUTP0970

OUTP0980

OUTP0990

OUTP1000

OUTP1010

OUTP1020

OUTP1030

OUTP1040

OUTP1050

OUTP1060

OUTP1070

OUTP1080

OUTP1090

OUTP1100

OUTP1110

OUTP1120

OUTP1130

OUTP1140

OUTP1150

OUTP1160

OUTP1170

OUTP1180

OUTP1190

OUTP1200

OUTP1210

OUTP1220

OUTP1230

OUTP1240

OUTP1250

OUTP1260

OUTP1270

OUTP1280

OUTP1290

OUTP1300

OUTP1310

OUTP1320

OUTP1330

98.

C	IF(PROB)7162,7162,7163	OUTP1350
7163	N3=0	OUTP1360
7162	WS=555.0	OUTP1370
	WRITE (N7)WS,CYCLE,N3	OUTP1380
	WRITE (N7)(Z(I),I=1,MZ)	OUTP1390
7131	WRITE (N7)(U(K),V(K),AMX(K),AIX(K),AIX(K),K=1,KMAXA)	OUTP1400
	GO TO 7140	OUTP1410
7140	CONTINUE	OUTP1420
	WRITE (N7)X(0),(X(K),TAU(K),K=1,IMAX)	OUTP1430
	WRITE (N7)(Y(K),K=0,JMAX)	OUTP1440
	WS=666.0	OUTP1450
		OUTP1460
		OUTP1470
C	WRITE PARTICLES ON DUMP TAPE FOR PIC RUN.	
C	IF(PROB)7150,7150,7161	OUTP1480
7150	DO 7160 I=1,N3	OUTP1490
	READ (N2)(AM(N),XL(N),YL(N),IW1(N),IW2(N),N=2,NPRI)	OUTP1500
	WRITE (N7)(AM(N),XL(N),YL(N),IW1(N),IW2(N),N=2,NPRI)	OUTP1510
7160	CONTINUE	OUTP1520
7161	WRITE (N7)WS,WS,WS	OUTP1530
	REWIND N7	OUTP1540
	WRITE (6,8120)T,NC	OUTP1550
	IWS=IMAX*JMAX+1	OUTP1560
	CALL SLITE (0)	OUTP1570
	DO 7517 I=1,IMAX	OUTP1580
	CALL SLITE (1)	OUTP1590
	J=JMAXA	OUTP1600
	K=IWS+I	OUTP1610
	DO 7517 JP=1,JMAX	OUTP1620
	J=J-1	OUTP1630
	K=K-IMAX	OUTP1640
7170	IF(AMX(K))9905,7517,7175	OUTP1650
7175	CALL SLITET(1,K000FX)	OUTP1660
	GO TO(7180,7185),K000FX	OUTP1670
C	PRINT OUT CELL QUANTITIES.	
7180	WRITE (6,8080)I,X(I),DX(I)	OUTP1680
7185	WRITE (6,8084)J,Y(J),DY(J),U(K),V(K), I,V(K),AMX(K)	U(K),AIX(K) OUTP1690
7517	CONTINUE	OUTP1700
	IF(Q000FL)7520,7520,7616	OUTP1710
7616	REWIND N2	OUTP1720
	GO TO 7520	OUTP1730
C	ERROR	OUTP1820
9901	NK=7013	OUTP1830
	GO TO 9999	OUTP1840
9902	NK=7014	OUTP1850
	GO TO 9999	OUTP1860
9904	NK=7005	OUTP1870
	GO TO 9999	OUTP1880
9905	Nk=7170	OUTP1890
		OUTP1900

99.

9999 WRITE (6,8888)NK,I,J,K,L,M,N
 PRINT 8888,NK,I,J,K,L,M,N
 CALL DUMP
 7520 RETURN

OUTP1910
 OUTP1920
 OUTP1930
 OUTP1940
 OUTP1950

C

FORMATS

80720FORMAT(1H ///6H THE =1PE16.9,7X,3HE =E16.9///5H M. =E11.5,5X,4HMX
 1 =E11.5,7X,7HM.+MX =E11.5)
 8073 FORMAT(1H0/17HOPARTICLES - - -112,4H DOT114,2H X114,6H TOTAL)
 80800FORMAT(1H0///3H0I=12,10X,2HX=1PE13.7,10X,3HDX=E13.7/3H0 J10X,1HY130
 1X,2H0Y12X,1HU13X,1HV12X,3HAID11X,3HAIX11X,3HAMD11X,3HAMX)
 8084 FORMAT(13,3X,1P8E14.7)
 8104 FORMAT(1H /31H THERE ARE NO MORE PACKAGES----)
 8120 FORMAT(1H ///18H TAPE DUMP AT TIMEF10.1,7X,5HCYCLEI4)
 8888 FORMAT(1H+/26H1OUTPUT ERROR IN STATEMENTI5,12X,12H INDICES ARE6I7)
 END

OUTP1960
 OUTP1970
 OUTP1980
 OUTP1990
 OUTP2000
 OUTP2010
 OUTP2060
 OUTP2070
 OUTP2080
 OUTP2090

5.2. FORTRAN IV LISTINGS OF OIL.

100.

*** NOTE, THE FOLLOWING SET OF DIMENSIONS, COMMON
AND EQUIVALENCE CARDS ARE TO BE USED FOR ALL
SUBROUTINES WITH THE EXCEPTION OF MAIN AND CARDS.

D I M E N S I O N

PH2 0020

PH2 0030

PH2 0040

PH2 0050

PH2 0060

PH2 0070

PH2 0080

PH2 0090

PH2 0100

PH2 0110

PH2 0120

PH2 0130

PH2 0140

PH2 0150

PH2 0160

PH2 0170

PH2 0180

PH2 0190

PH2 0200

PH2 0210

PH2 0220

PH2 0230

PH2 0240

PH2 0250

PH2 0260

PH2 0270

PH2 0280

PH2 0290

PH2 0300

PH2 0390

PH2 0400

PH2 0410

PH2 0420

PH2 0430

PH2 0440

PH2 0450

PH2 0460

PH2 0470

PH2 0480

PH2 0490

PH2 0500

PH2 0510

PH2 0520

```

DIMENSION AM(130), XL(130), YL(130),
1U(3500),V(3500),AMX(3500),AIX(3500),
2P(3500),
3 IW1(130), IW2(130),
4DX(52), X(53), XX(54), DY(100), Y(100), YY(101),
5TAB(15), AMK(15), PK(15), QK(15), Z(150), IZ(150),
6TAU(52), PL(200), PR(200), UL(200), UR(200),
7FLEFT(100),YAMC(100), SIGC(100), GAMC(100)
COMMON Z ,XX ,UR ,PR ,YY
COMMON AID ,AIX ,AM ,AMD ,AMX ,AREA
COMMON BIG ,BOUNCE ,DDXN ,DDVK ,DKE ,DVK
COMMON DX ,DY ,E ,FD ,FS ,FX
COMMON OUT ,P ,PABOVE ,PBLO ,PIDTS ,PPABOV
COMMON PRR ,PUL ,QDT ,RC ,REZ ,RHO
COMMON RL,RR,SIG,Q000FL,SWITCH ,TABLM
COMMON TAU ,TAUDTS ,TAUDTX ,U ,UK ,URR
COMMON UT ,UU ,UUU ,UTEF ,UVMAX ,V
COMMON VABOVE ,VBLO ,VEL ,VK ,VT ,VTEF
COMMON VV ,VVABOV ,VVBLO ,W2 ,W3 ,WPS
COMMON WS ,WSA ,WSB ,WSC ,XL ,XLF
COMMON XN ,XR ,YL ,YLN ,YN ,YU
COMMON ZMAX ,I ,II ,IN ,IR ,IWS
COMMON IWSA ,IWSB ,IWSC ,IW1 ,J ,JN
COMMON JP ,JR ,K ,KN ,KP ,KR
COMMON KRM ,L ,M ,MA ,MB ,MC
COMMON MD ,ME ,MZ ,N ,NK ,NKMAX
COMMON NK1 ,NO ,NR ,IW2

```

E Q U I V A L E N C E

```

OEQUIVALENCE (Z,IZ,PROB), (Z(2),CYCLE), (Z(3),DT),
1(Z(4),PRINTS), (Z(5),PRINTL), (Z(6),DUMPT7), (Z(7),CSTOP),
2(Z(8),PIDY), (Z(9),TMZ), (Z(10),GAM), (Z(11),GAMD),
3(Z(12),GAMX), (Z(13),ETH), (Z(14),FFA), (Z(15),FFB),
4(Z(16),TMDZ), (Z(17),TMXZ), (Z(18),XMAX), (Z(19),TXMAX),
5(Z(20),TYMAX), (Z(21),AMDM), (Z(22),AMXM), (Z(23),DNN),
6(Z(24),DMIN), (Z(25),FEF), (Z(26),DTNA), (Z(27),CVIS),
7(Z(28),NPR), (Z(29),NPRI), (Z(30),NC), (Z(31),NPC),
8(Z(32),NRC), (Z(33),IMAX), (Z(34),IMAXA), (Z(35),JMAX),
9(Z(36),JMAXA), (Z(37),KMAX), (Z(38),KMAXA), (Z(39),NMAX)

```

OEQUIVALENCE	(Z(40),ND),	(Z(41),KDT),	(Z(42),IXMAX),	PH2 0530
1(Z(43),NOJ),	(Z(44),NOPR),	(Z(45),NIMAX),	(Z(46),NJMAX),	PH2 0540
2(Z(47),I1),	(Z(48),I2),	(Z(49),I3),	(Z(50),I4),	PH2 0550
3(Z(51),N1),	(Z(52),N2),	(Z(53),N3),	(Z(54),N4),	PH2 0560
4(Z(55),N5),	(Z(56),N6),	(Z(57),N7),	(Z(58),N8),	PH2 0570
5(Z(59),N9),	(Z(60),N10),	(Z(61),N11),	(Z(62),NRM),	PH2 0580
6(Z(63),TRAD),	(Z(64),XNRG),	(Z(65),SN),	(Z(66),DXN),	PH2 0590
7(Z(67),RADER),	(Z(68),RADET),	(Z(69),RADER),	(Z(70),DTRAD),	PH2 0600
8(Z(71),REZFCT),	(Z(72),RSTOP),	(Z(73),SHELL),	(Z(74),SBOUND),	PH2 0610
9(Z(75),TOZONE),	(Z(76),ECK),	(Z(77),SBOUND),	(Z(78),X1),	PH2 0620
OEQUIVALENCE	(Z(79),X2),	(Z(80),Y1),	(Z(81),Y2),	PH2 0630
1(Z(82),CABL),	(Z(83),VISC),	(Z(84),T),	(Z(85),GMAX),	PH2 0640
2(Z(86),WSGD),	(Z(87),WSGX),	(Z(88),GMADR),	(Z(89),GMAXR),	PH2 0650
3(Z(90),S1),	(Z(91),S2),	(Z(92),S3),	(Z(93),S4),	PH2 0660
4(Z(94),S5),	(Z(95),S6),	(Z(96),S7),	(Z(97),S8),	PH2 0670
5(Z(98),S9),	(Z(99),S10)			PH2 0680
				PH2 0690
OEQUIVALENCE	(XX(2),X(1)),	(UR,UL,FLEFT),	(UR(100),YAMC),	PH2 0700
1(PR(100),SIGC),	(PR,PL,GAMC),	(UR,TAB),		PH2 0710
2(UR(16),AMK),	(UR(31),PK),	(UR(46),QK),	(YY(2),Y(1))	PH2 0720
				PH2 0730
				PH2 0740

NOTE, THERE ARE 2 SPECIAL SUBROUTINES(FORTRAN 4) USED
IN THE OIL CODE, SUBROUTINE SLITE SERVES THE SAME FUNCTION
AS TURNING ON SENSE LIGHTS, AND SLITET SERVES THE
FUNCTION OF TESTING THE SENSE LIGHTS.

NOTE, IF AN ERROR(SEE THE END OF THE SUBROUTINES) OCCURS, THE
SUBROUTINE WILL CALL FOR A DUMP. BY CHECKING THE
VALUES OF NR AND NK , ONE CAN READILY IDENTIFY THE
STATEMENT NUMBER AND THE SUBROUTINE WHERE THE ERROR OCCURRED.

NK WILL CONTAIN THE STATEMENT NUMBER, AND NR IS
AN IDENTIFICATION FOR THE SUBROUTINE AS FOLLOWS,

NR=1	INPUT
NR=2	CDT
NR=3	PH1
NR=4	PH2
NR=6	EDIT

102.

\$IBFTC MAIN LIST,DECK,REF

CMAIN

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***** NOTE 1 MATERIAL ONLY (X1) *****

INPUT READS OIL DUMP TAPE OR
WILL CALL SUBROUTINE SET*UP WHICH
WILL MAKE A DUMP TAPE FOR CERTAIN TYPES OF PROBLEM
(SEE SECTION ON SET*UP)
ALSO CALCULATES DX AND DY AND EQUATION OF STATE DATA
CALL INPUT

CDT ROUTINE CALCULATES DT(HYDRO TIME STEP)
AND PRESSURES, ADVANCE CYCLE NO. ETC.

10 CALL CDT

IN EDIT, DETERMINE WHETHER TO EXECUTE A LONG
PRINT, A SHORT PRINT, A TAPE DUMP, ETC. AND
CALCULATE TOTAL ENERGY IN SYSTEM(COMPARE
WITH ETH) TOTAL MASS, INTEGRATE TOTAL
COMPONENTS OF MOMENTA.

CALL EDIT

CALL SLITET(1,K000FX)

SENSE LITE 1 SIGNIFIES THIS
IS THE LAST CYCLE OF THIS RUN \$\$\$\$\$\$\$\$\$\$\$\$\$\$\$\$
LITE TURNED ON IN THE EDIT ROUTINE *****

GO TO(30,20),K000FX

PH1, INTEGRATE THE MOMENTA EQS. INTEGRATE
ENERGY EQUATION(ONLY CHANGES DUE TO WORK
TERMS). NO MOVEMENT OF MASS HERE

20 CALL PH1

TRANSPORT MASS ACROSS BOUNDARIES (SOLVE
MASS TRANSPORT EQ.) TRANSPORT TERMS IN
THE MOMENTA AND ENERGY EQS. LEFT OUT OF
PH1, HERE APPROXIMATED BY MASS MOVEMENT. CONSERVE
MASS, MOMENTA AND TOTAL ENERGY.

CALL PH2

GO TO 10

30 CALL EXIT

END

MAIN0010

MAIN0020

MAIN0030

MAIN0050

MAIN0060

MAIN0070

MAIN0080

MAIN0090

MAIN0100

MAIN0110

MAIN0120

MAIN0130

MAIN0140

MAIN0150

MAIN0160

MAIN0170

\$IBFTC INPUT LIST,DECK,REF	
SUBROUTINE INPUT	
C	INPU0010
C	INPU0760
C	INPU0900
C TURN ON SENSE LITE 3.	
CALL SLITE (3)	INPU0980
C	INPU0990
C READ HEADER CARD (COLUMNS 2-72).	
READ (5,8004)IWS	INPU1000
WRITE (6,8004)IWS	INPU1010
C CALL DATA.	
6 CALL CARDS	INPU1020
C IF PK(3) = OR GREATER THAN ZERO, CALL ROUTINE	
C SET-UP, OTHERWISE, BINARY OIL TAPE HAS BEEN MADE.	
C READ IN DATA FROM OIL DUMP TAPE, OR	
C GENERATE A DUMP TAPE FOR OIL, AND	
C CALCULATE DX AND DY FROM THE X AND	
C Y VALUES FROM TAPE.	
IF(PK(3))8687,8888,8888	INPU1030
8888 CALL CARDS	INPU1040
CALL SETUP	INPU1050
8887 CONTINUE	INPU1060
C	INPU1070
C READ TAPE	INPU1080
C GO READ BINARY TAPE.	
GO TO 1000	INPU1090
C	INPU1100
C READ IN REMAINING INPUT CARDS	INPU1110
10 CONTINUE	INPU1120
CALL CARDS	INPU1130
GO TO 2000	INPU1140
C	INPU1150
C SET THE PRESSURES TO ZERO.	
40 DO 45 K=1,KMAXA	INPU1160
45 P(K)=0.0	INPU1170
C INTEGRATE BACKWARDS ON CYCLE, TIME AND NO. OF	
C CYCLES BETWEEN ENERGY CHECK, SINCE THESE	
C ARE ADVANCED IN CDT.	
T=T-DTNA	INPU1180
NC=NC-1	INPU1190
CYCLE=NC	INPU1200
NPC=NPC-1	INPU1210
UVMAX=0.0	INPU1220
C CALCULATE THE DX'S, SINCE THESE ARE NOT ON	
C TAPE.	
DO 50 I=1,IMAX	INPU1230
50 DX(I)=X(I)-X(I-1)	INPU1240
C CALCULATE THE DY'S, SINCE THESE ARE NOT ON	
C TAPE.	
DO 55 J=1,JMAX	INPU1250

104.

55	DY(J)=Y(J)-Y(J-1)	INPU1260
	J=MZ-8	
C	PRINT Z BLOCK.	
62	DO 80 I=1,J,8	
	K=I+7	INPU1290
	DO 65 J=I,K	INPU1300
	IF(Z(J))70,65,70	INPU1310
65	CONTINUE	INPU1320
	GO TO 80	INPU1330
70	K=I+7	INPU1340
	WRITE (6,8111)I,(Z(L),L=I,K)	INPU1350
80	CONTINUE	INPU1360
	GO TO 10000	INPU1370
C		INPU1380
C		INPU1390
C		INPU1400
C	READ BINARY TAPE.	
1000	MZ=150	INPU1410
	IWS=0	INPU1420
1003	REWIND 7	
1004	READ(7)PR(1),PR(2),N3	
	NR=N3+5	INPU1450
1006	IF(PR(1)-555.0)1010,1016,1010	INPU1460
1010	IWS=IWS+1	INPU1470
1011	IF(MOD(IWS,3))9902,9902,1003	INPU1480
1016	IF(PR(2))1010,1018,1018	INPU1490
C	CHECK HERE FOR THE CORRECT CYCLE NUMBER.	
1018	IF(PK(2)-PR(2))1023,1023,1020	INPU1500
C	SKIP OVER, LOOK AT NEXT CYCLE.	
1020	DO 1022 L=2,NR	INPU1510
1022	READ(7)	
	GO TO 1004	INPU1530
1023	READ(7)(Z(I),I=1,MZ)	
C	CHECK FOR THE CORRECT PROBLEM NO.	
	IF(ABS(PROB-PK(1))-.01)1024,1024,9901	INPU1550
1024	READ(7)(U(I),V(I),AMX(I),AIX(I),P(I),I=1,KMAX)	
	READ(7)X(0),X(I),TAU(I),I=1,IMAX)	
	READ(7)(Y(I),I=0,JMAX)	
C	NOTE, INITIALIZE N1 TO 2, AND N2 TO 3	
C	NOTE, N1 AND N2 ARE ONLY USED FOR BOOK-KEEPING *****	
	N1=2	
	N2=3	
C	NOTE, IF PROBLEM NO. IS NEGATIVE, THIS IS	
C	A PIC TRANSPORT, CHECK YOUR PH2 ROUTINE,	
C	AND READ THE PARTICLES FROM TAPE ONTO	
C	SCRATCH TAPE N1.	
	IF(PROB)1,1,1034	INPU1590
1	REWIND 2	
	REWIND 3	
	DO 1025 I=1,N3	INPU1620

```

      READ(7)(AM(N),XL(N),YL(N),IW1(N),IW2(N),N=2,N4)
      WRITE(2)(AM(N),XL(N),YL(N),IW1(N),IW2(N),N=2,N4)
1025 CONTINUE
1034 READ(7)PR(1),PR(2),PR(3)
      REWIND 2
1036 IF(PR(1)-555.0)9904,1040,1038
1038 IF(PR(2)-666.0)9905,1040,9905
1040 GO TO 10
C**** END OF READ TAPE *****
C
C
C      CALCULATE MAX. GAMMA AND GAMMA/(GAMMA-1.).
C
2000 IF(WSGX)9906,2010,2005
2005 GAMX=1.0/(WSGX-1.0)
2010 WSGX=(GAMX+1.0)/GAMX
      GMAXR=GAMX*WSGX
2012 IF(WSGD)9907,2020,2015
2015 GAMD=1.0/(WSGD-1.0)
2020 WSGD=(GAMD+1.0)/GAMD
      GMADR=GAMD*WSGD
      GMAX=WSGD
      IF(WSGD-WSGX)2025,2030,2030
2025 GMAX=WSGX
2030 GO TO 40
C**** END OF R E S *****
C
C
C      ERROR
9901 NK=1023
      GO TO 9999
9902 NK=1011
      GO TO 9999
9904 NK=1036
      GO TO 9999
9905 NK=1038
      GO TO 9999
9906 NK=2000
      GO TO 9999
9907 NK=2012
9999 NR=1
      CALL DUMP
C
10000 RETURN
C
C      FORMATS
8000 FORMAT(7E10.3,I2)
80040FORMAT(I1,7I1H
1
8111 FORMAT(I4,8O14)
END

```

INPU1650

INPU1680

INPU1690

INPU1700

INPU1710

INPU1720

INPU1730

INPU1740

INPU1750

INPU1760

INPU1770

INPU1780

INPU1790

INPU1800

INPU1810

INPU1820

INPU1830

INPU1840

INPU1850

INPU1860

INPU1870

INPU1880

INPU1890

INPU1900

INPU1910

INPU1920

INPU1930

INPU1940

INPU1950

INPU1960

INPU1970

INPU1980

INPU1990

INPU2000

INPU2010

INPU2020

INPU2030

INPU2040

INPU2050

INPU2060

INPU2070

INPU2080

INPU2090

INPU2100

INPU2110

\$IBFTC CARDS LIST,DECK,REF

SUBROUTINE CARDS

DIMENSION TABLE(1),CARD(7),LABEL(1)

COMMON TABLE

C A 2 IN COLUMN 1, ROUTINE WILL FIX THE
C FLOATING PT. NO.
C A 1 IN COLUMN 1, MEANS THIS IS LAST CARD TO
C READ IN.

EQUIVALENCE(TABLE(1),LABEL(1))

WRITE (6,10)

1 READ (5,11)IEND,LOC,NUMWPC,(CARD(I),I=1,NUMWPC)

WRITE (6,12)IEND,LOC,NUMWPC,(CARD(I),I=1,NUMWPC)

DO 4 I=1,NUMWPC

J=LOC+I-1

IF(IEND-2)2,5,2

5 LABEL(J)=IFIX(CARD(I))

GO TO 4

2 TABLE(J)=CARD(I)

4 CONTINUE

IF(IEND-1)1,3,1

3 RETURN

C FORMATS

10 FORMAT(20H1 OIL INPUT CARDS///)

11 FORMAT(11,15,11,OP7E9.4)

12 FORMAT(1H 14,17,13,1P7E14.6)

END

CARD0010

CARD0020

CARD0030

CARD0050

CARD0070

CARD0080

CARD0090

CARD0100

CARD0110

CARD0120

CARD0130

CARD0140

CARD0150

CARD0160

CARD0170

CARD0180

CARD0190

CARD0210

CARD0220

CARD0230

\$IBFTC SETUP LIST,DECK,REF	
SUBROUTINE SETUP	SETU0010
C WILL ONLY GENERATE (1) MATERIAL.	
C PACKAGES MUST BE RECTANGLES.	
C ASSUMPTION OF = DX AND = DY	SETU0980
C LOAD PK(4)=1.	
M=PK(4)	SETU0990
C LOAD PK(5)=RIGHT BOUNDARY OF PELLET(I).	
MA=PK(5)	SETU1000
C LOAD PK(6)=BOTTOM(J)+1 OF PELLET.	
MB=PK(6)	SETU1010
C LOAD PK(7)=TOP(J) OF PELLET.	
MC=PK(7)	SETU1020
C LOAD PK(8)=1.	
MD=PK(8)	SETU1030
C LOAD PK(9)=RIGHT(I) BOUNDARY OF TARGET.	
ME=PK(9)	SETU1040
C LOAD PK(10)=BOTTOM(J)+1 OF TARGET.	
MZ=PK(10)	SETU1050
C LOAD PK(11)=TOP(J) OF TARGET.	
N=PK(11)	SETU1060
C LOAD INITIAL DENSITY INTO Z(111).	
RHO=Z(111)	SETU1070
C LOAD INITIAL PELLET VELOCITY INTO Z(112).	
VTEF=Z(112)	SETU1080
KMAX=IMAX*JMAX+1	SETU1090
KMAXA=XMAX+1	SETU1100
JMAXA=JMAX+1	SETU1110
IMAXA=IMAX+1	SETU1120
C CLEAR ALL CELL ARRAYS.	
DO 1 K=1,KMAX	SETU1130
U(K)=0.0	SETU1140
V(K)=0.0	SETU1150
P(K)=0.0	SETU1160
AMX(K)=0.0	SETU1170
AIX(K)=0.0	SETU1180
1 CONTINUE	SETU1190
DX(1)=DX(1)	SETU1200
X(1)=DX(1)	SETU1210
WS=X(1)**2	SETU1220
PIDY=3.1415927	SETU1230
TAU(1)=WS*PIDY	SETU1240
C CALCULATE DX,X,TAU	
DO 10 I=2,IMAX	SETU1250
X(I)=X(I-1)+DX(1)	SETU1260
DX(I)=DX(1)	SETU1270
WSA=X(I)**2	SETU1280
TAU(I)=PIDY*(WSA-WS)	SETU1290
WS=WSA	SETU1300
10 CONTINUE	SETU1310

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	Y(1)=DY(1)	SETU1320
C	CALCULATE DY AND Y.	
	DO 20 J=2,JMAX	SETU1330
	Y(J)=Y(J-1)+DY(1)	SETU1340
	DY(J)=DY(1)	SETU1350
20	CONTINUE	SETU1360
	ETH=0.0	SETU1370
	DO 30 I=M,MA	SETU1380
	K=(MB-1)*IMAX+I+1	SETU1390
C	CALCULATE MASS, AND VELOCITY OF PELLET.	
	DO 40 J=MB,MC	SETU1400
	AMX(K)=RHO*DY(J)*TAU(I)	SETU1410
	V(K)=VTEF	SETU1420
C	CALCULATE TOTAL ENERGY (ETH.)	
	ETH=ETH+AMX(K)*(V(K)**2)/2.0	SETU1430
40	K=K+IMAX	SETU1440
30	CONTINUE	SETU1450
C	CALCULATE MASS OF TARGET.	
	DO 50 I=MD,ME	SETU1460
	K=(MZ-1)*IMAX+I+1	SETU1470
	DO 60 J=MZ,N	SETU1480
	AMX(K)=RHO*DY(J)*TAU(I)	SETU1490
60	K=K+IMAX	SETU1500
50	CONTINUE	SETU1510
	IMAX=IMAX	SETU1520
	JMAX=JMAX	SETU1530
	SHELL=2.0	SETU1540
	CYCLE=0.0	SETU1550
	DT=0.0	SETU1560
	NMAX=0	SETU1570
	N1=2	SETU1580
	N2=3	SETU1590
	N3=0	SETU1600
	N4=127	SETU1610
	XMAX=X(IMAX)	SETU1620
	TXMAX=XMAX*2.0	SETU1630
	YMAX=Y(JMAX)	SETU1640
	TYMAX=YMAX*2.0	SETU1650
	REWIND 7	
	WS=555.0	SETU1670
C	WRITE OUTPUT FOR OIL ON TAPE.	
	WRITE (7)WS,CYCLE,N3	
	WRITE (7){Z(I),I=1,150}	
	WRITE (7){U(I),V(I),AMX(I),AIX(I),P(I),I=1,KMAXA}	
	WRITE (7){X(I),Y(I),TAU(I),I=1,IMAX}	
	WRITE (7){Y(I),I=0,JMAX}	
	WS=666.0	SETU1730
	WRITE (7)WS,WS,WS	
	REWIND 7	
	RETURN	SETU1760
	END	

\$18FTC CDT LIST,DECK,REF
SUBROUTINE CDT

CDT 0010
CDT 0020
CDT 0990
CDT 1000
CDT 1010
CDT 1020

C
C
C =====
C
C
C CHECK COURANT CONDITION AND PARTICLE
C VELOCITY.
C RECORD I AND J OF ZONE WHERE DT IS BEING
C CONTROLLED.
3000 VEL=0.0
3005 DO 3050 I=1,I1
3010 K=I+1
3015 DO 3050 J=1,I2
I=I
J=J
3020 IF(AMX(K))9901,3050,3025
C
C CALCULATE PRESSURES FROM EQUATION OF STATE(ES).
3025 CALL ES
C
3030 IF(ABS(P(K))-1.0E-20)3035,3035,3040
3035 P(K)=0.0
3040 IF(WSGX-VEL)3050,3050,3045
3045 VEL=WSGX
3050 K=K+IMAX
3055 KDT=1
UVMAX=-1.0
3070 DO 3255 I=1,I1
3075 K=I+1
3095 DO 3255 J=1,I2
3100 KP=K+IMAX
IF(AMX(K))9901,3255,4
C IF RHO(K) IS LESS THAN Z(138), CELL K
C WILL BE BYPASSED FOR STABILITY CHECK.
4 IF(AMX(K)/(TAU(I)*DY(J))-Z(138))3255,3255,3115
3115 SIG=DX(I)
3120 IF (DY(J)-SIG)3125,3130,3130
3125 SIG=DY(J)
C C=SPEED OF SOUND FOR POLYTROPIC GAS AS
C THE SQ. ROOT OF (GAMMA*P/RHO).
C HERE CALCULATE THE SPEED OF SOUND FOR
C THE EQUATION OF STATE
C AS THE SQ. ROOT OF DP/DRHO.
3130 IF(Z(148))4000,4000,4001
4000 WS=SQRT(GMAX*TAU(I)*DY(J)*ABS(P(K))/(AMX(K)))
GO TO 3205
4001 WSA=ABS(P(K))*1.E+4
WS=Z(148)+Z(149)*(WSA**Z(150))

CDT 1030
CDT 1040
CDT 1050
CDT 1060
CDT 1070
CDT 1080
CDT 1090
CDT 1100
CDT 1110
CDT 1120
CDT 1130
CDT 1140
CDT 1150
CDT 1160
CDT 1170
CDT 1180
CDT 1190
CDT 1200
CDT 1210
CDT 1220
CDT 1230
CDT 1240
CDT 1250
CDT 1260
CDT 1270
CDT 1280
CDT 1290
CDT 1300
CDT 1310
CDT 1320
CDT 1330

110.

WS=WS*1.E-3	CDT 1340
3205 WS=WS/SIG	CDT 1350
3210 IF(UVMAX-WS)3215,3220,3220	CDT 1360
3215 N10=I	CDT 1370
N11=J	CDT 1380
UVMAX=WS	CDT 1390
3220 IF(NMAX)1,1,2	CDT 1400
C EULERIAN CHECK FOR RADIAL PARTICLE VELOCITY.	
1 CONTINUE	CDT 1410
3 WS=ABS(U(K))/TAU(I)*X(I)/.5*PI*DY	CDT 1420
GO TO 3225	CDT 1430
C PIC CHECK FOR RADIAL PARTICLE VELOCITY.	
2 WS=ABS(U(K))/DX(I)	CDT 1440
3225 IF(UVMAX-WS)3230,3235,3235	CDT 1450
3230 UVMAX=WS	CDT 1460
N10=I	CDT 1470
N11=J	CDT 1480
3235 WS=ABS(V(K))/DY(J)	CDT 1490
3240 IF(UVMAX-WS)3245,3250,3250	CDT 1500
3245 N10=I	CDT 1510
N11=J	CDT 1520
UVMAX=WS	CDT 1530
3250 CONTINUE	CDT 1540
3255 K=K+IMAX	CDT 1550
IF(UVMAX)9912,9912,3260	
C FOR OPTIONS ON CABLN, CHECK	
C SECTION 3.4 IN GAMD-5580.	
3260 IF(CABLN)90,91,3300	CDT 1560
90 DT=.5/VEL/UVMAX*Z(139)	CDT 1570
GO TO 3295	CDT 1580
91 WS=UVMAX*DT	CDT 1590
WSA=0.5/VEL	CDT 1600
3265 IF(FFA-WSA)3270,3276,3270	CDT 1610
3270 FFA=WSA	CDT 1620
3276 IF(WS-FFA)3285,3300,3280	CDT 1630
3280 DT=DT/WS*FFB/0.9	CDT 1640
GO TO 3295	CDT 1650
3285 IF(WS-FFB)3290,3290,3300	CDT 1660
3290 DT=DT*FFA/WS*0.9	CDT 1670
3295 KDT=0	CDT 1680
C INTEGRATE THE TIME AND CYCLE COUNTER.	
3300 T=T+DTNA	CDT 1690
85 IF(DTRAD)9911,80,81	CDT 1700
80 NR=NRN	CDT 1710
84 WS=NR	CDT 1720
TRAD=DT/WS	CDT 1730
GO TO 82	CDT 1740
81 IWS=DT/DTRAD	CDT 1750
NR=IWS+1	CDT 1760
83 IF(NR-NRN)84,84,80	CDT 1770

82 NC=NC+1		CDT 1780
CYCLE=NC		CDT 1790
NPC=NPC+1		CDT 1800
3305 IF(T)9909,3320,3310		CDT 1810
3310 IF(KDT)9910,3315,3320		CDT 1820
3315 WRITE (6,8000)T,DTNA,DT		CDT 1830
3320 DTNA=DT		CDT 1840
GO TO 3325		CDT 1850
C	NEGATIVE MASS	CDT 1860
9901 NK=3020		CDT 1870
GO TO 9999		CDT 1880
9909 NK=3305		CDT 1890
GO TO 9999		CDT 1900
9910 NK=3310		CDT 1910
GO TO 9999		CDT 1920
C	THE DT WILL BE 0. OR NEGATIVE ,STOP	
9912 NK=1		
GO TO 9999		
9911 NK=85		
9999 NR=2		CDT 1930
CALL DUMP		CDT 1940
3325 RETURN		CDT 1950
80000FORMAT (17HOCHANGE DT ... T=1PE9.3,11H	DT(N)=1PE9.3,13H	CDT 1960
1(N+1)=1PE9.3)		DTCDT 1970
END		CDT 1980
		CDT 1990

112.

```
$IBFTC PH1      LIST,DECK,REF
SUBROUTINE PH1
```

PH1 0010
PH1 0900

C VELOCITIES, ENERGIES, PRESSURES ARE AT THE
C CENTER OF THE CELL.
C (2) PASSES THRU PH1 ARE REQUIRED. NO
C MASS IS MOVED IN PH1.
C ***** NOTE 1 MATERIAL ONLY (X) *****

PH1 0980
PH1 0990
PH1 1000
PH1 1010
PH1 1020
PH1 1030
PH1 1040
PH1 1050
PH1 1060
PH1 1070
PH1 1080

C
 C
 C
 C
 C

```

NRT=0
NRC=0
UU=1.E+15
UT=0.0
C YOU WILL GET BACK HERE IF AIX WAS LESS
C THAN 0. AND PROVIDED SN=0.

```

PHI 1090

```

8000 VEL=1.0
C      INITIALIZE MID-POINTS OF FIRST AND SECOND
C      CELL IN R DIRECTION.

```

```
3301 RC=DX(1)/2.0
      RR=(X(1)+X(2))/2.0
```

PH1 1100
PH1 1110
PH1 1120

```

3304 K=2
C  AXIS OF SYMMETRY BOUNDARY CONDITIONS.
DO 3302 J=1,JMAX
PL(J)=P(K)
UL(J)=0.0

```

PH1 1130
PH1 1140
PH1 1150
PH1 1160

```

3302 K=K+IMAX
C   FIRST PASS THRU, CALCULATE U AND V AT
C   CYCLE N+1, AND THE WORK TERMS USING U AND V
C   FROM CYCLE N.
C   SECOND PASS THRU, CALCULATE ONLY THE
C   CONTRIBUTION TO THE CHANGE IN INTERNAL ENERGY
C   FROM WORK TERMS EVALUATED FROM U AND V
C   AT CYCLE N+1.

```

PH1 1170
PH1 1180
PH1 1190

```

DO 3360 I=1,I1
K=I+1
IF(CVIS)7002,7003,7003
C   BOTTOM BOUNDARY IS TRANSMISSIVE.

```

```

7002 VBLO=V(K)
      PBLO=0.0
      GO TO 7004

```

PH1 1200
PH1 1210
PH1 1220

C BOTTOM BOUNDARY IS REFLECTIVE.

```

7003 VBLO=0.0
      PBLO=P(K)

```

PH1 1230
PH1 1240
PH1 1250

```

7004 TAUDTS=TAU(I)*DT
C      I1= MAX.(I) OF DISTURBANCE IN R DIRECTION.

```

C	I2= MAX(J) OF DISTURBANCE IN Z DIRECTION.	
C	DO LOOP IN J DIRECTION	
	DO 3348 J=1,I2	PH1 1260
	PIDTS=1.0/(PIDY*DT*DY(J))	PH1 1270
C	K= INDEX OF CELL IN QUESTION.	
C	N= INDEX OF CELL ABOVE.	
	N=K+IMAX	PH1 1280
	3305 IF(AMX(K))9902,3340,3306	PH1 1290
	3306 IF(IMAX-I)9903,3311,3310	PH1 1300
	3310 IF(AMX(K+1))9904,3312,3314	PH1 1310
C	WE ARE AT THE RIGHT BOUNDARY, SET PRESSURE	
C	GRADIENT TO 0. IN R DIRECTION, MODIFY ETH.	
C	FOR RIGHT BOUNDARY BEING TRANSMITTIVE.	
	3311 PRR=PL(J)	PH1 1320
	3307 ETH=ETH-PRR*U(K)/PIDTS*RC	PH1 1330
	GO TO 3313	PH1 1340
C	RIGHT BOUNDARY CONDITION FOR THE MOMENTUM EQ.	
C	ADJACENT TO EMPTY CELL.	
	3312 PRR=0.0	PH1 1350
	3313 URR=RC*U(K)	PH1 1360
	GO TO 3316	PH1 1370
C	CALCULATE PRESSURE AT INTERFACE(I) AND	
C	(RU) FOR WORK TERM.	
	3314 PRR=(P(K)+P(K+1))/2.0	PH1 1380
	3315 URR=(U(K)*RC+U(K+1)*RR)/2.0	PH1 1390
	3316 IF(JMAX-J)9905,3318,3320	PH1 1400
C	SET PRESSURE GRADIENT TO 0. THIS IS FOR TOP	
C	BOUNDARY BEING TRANSMITTIVE.	
	3318 PABOVE=PBLO	PH1 1410
C	MODIFY ETH FOR TOP BOUNDARY CONDITION.	
	3319 ETH=ETH-PABOVE*V(K)/2.0*TAUDTS	PH1 1420
	GO TO 3323	PH1 1430
	3320 IF(AMX(N))9906,3322,3324	PH1 1440
C	TOP BOUNDARY CONDITION (EMPTY CELL ABOVE.)	
C	TOP BOUNDARY CONDITION FOR VELOCITY (EMPTY CELL ABOVE).	
	3322 PABOVE=0.0	PH1 1450
	3323 VABOVE=V(K)	PH1 1460
	GO TO 3328	PH1 1470
C	CALCULATE PRESSURE AT INTERFACE(J)	
	3324 PABOVE=(P(K)+P(N))/2.0	PH1 1480
	IF(CVIS)7001,3325,3325	PH1 1490
	7001 IF(1-J)3325,7000,9905	PH1 1500
C	BOTTOM BOUNDARY IS TRANSMITTIVE, SET PRESSURE	
C	GRADIENT TO 0.	
C	AND MODIFY ETH.	
	7000 PBLO=PABOVE	PH1 1510
	ETH=ETH+PBLO*V(K)/2.0*TAUDTS	PH1 1520
C	VELOCITY AT INTERFACE(J)	
	3325 VABOVE=(V(K)+V(N))/2.0	PH1 1530
	3328 IF(VEL)9907,3404,3400	PH1 1540

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C	COMPUTE DELTA U AND DELTA V.	
3400	V(K)=V(K)+(P8LO-PABOVE)*TAUDTS/(AMX(K))	PH1 1550
	IF(ABS(V(K))-1.E-08)3401,3401,3402	PH1 1560
3401	V(K)=0.0	PH1 1570
3402	U(K)=U(K)+(PL(J)-PRR)/(AMX(K))*RC/PIDTS*2.0	PH1 1580
	IF(ABS(U(K))-1.E-08)3403,3403,3404	PH1 1590
3403	U(K)=0.0	PH1 1600
C	CHECK FOR ADVANCING COUNTERS OF THE ACTIVE	
C	GRID IN THE R DIRECTION.	
3404	IF(I-I1)6016,6005,6005	PH1 1610
6005	IF(U(K))6605,6606,6605	PH1 1620
6605	NRC=1	PH1 1630
6606	IF(V(K))6607,6004,6607	PH1 1640
6607	NRC=1	PH1 1650
6004	IF(AIX(K))6015,6016,6015	PH1 1660
6015	NRC=1	PH1 1670
6016	CONTINUE	PH1 1680
C	HERE CALCULATE CHANGE IN INTERNAL ENERGY	
C	DUE TO WORK TERMS ONLY.	
	HS=(VBLO-VABOVE)*TAUDTS/2.0*P(K)	PH1 1690
	RHO=WS+(UL(J)-URR)/PIDTS*P(K)	PH1 1700
C	CONVERT TO SPECIFIC INTERNAL ENERGY.	
3332	WSX=AIX(K)*RHO/AMX(K)	PH1 1710
	GO TO 1000	PH1 1720
C	CHECK FOR NEGATIVE INTERNAL ENERGIES.	
1000	IF(WSX)1011,1001,1001	PH1 1730
1001	AIX(K)=WSX	PH1 1740
	GO TO 3342	PH1 1750
1011	UT=1.0	PH1 1760
C	COMPUTE NEW DT(STORE IN UU) ASSUMING	
C	THAT DI/DT WILL BE THE SAME FOR A SMALLER	
C	TIME STEP, THE NEW DT IS CHOSEN SUCH	
C	THAT AIX(AT N+1)=2/3 OF AIX(N).	
	WSA=2.0*AIX(K)/3.0*DT/(AIX(K)-WSX)	PH1 1770
1013	IF(WSA-UU)1014,1001,1001	PH1 1780
1014	UU=WSA	PH1 1790
	GO TO 1001	PH1 1800
C	CELL (K) IS EMPTY, SET INTERFACE QUANTITIES,	
C	ASSUMING CELL TO THE RIGHT AND TOP ARE	
C	NOT VOID.	
3340	PRR=0.0	PH1 1810
	URR=U(K+1)*RR	PH1 1820
	PABOVE=0.0	PH1 1830
	VABOVE=V(N)	PH1 1840
C	SET RIGHT QUANTITIES TO THE LEFT (FOR NEXT	
C	COLUMN SWEEP) AND SET ABOVE QUANTITIES TO	
C	BELOW FOR NEXT CELL ABOVE.	
3342	VSLO=VABOVE	PH1 1850
	PL(J)=PRR	PH1 1860
	UL(J)=URR	PH1 1870

115.

K=N	PH1 1880
3348 PBLO=PABOVE	PH1 1890
LL=K-IMAX	PH1 1900
C CHECK FOR ADVANCING COUNTERS OF THE ACTIVE	
C GRID IN Z DIRECTION.	
IF(U(LL))6000,6001,6000	PH1 1910
6000 NRT=1	PH1 1920
6001 IF(V(LL))6002,6003,6002	PH1 1930
6002 NRT=1	PH1 1940
6003 IF(AIX(LL))6017,6018,6017	PH1 1950
6017 NRT=1	PH1 1960
6018 CONTINUE	PH1 1970
3355 RC=RR	PH1 1980
RR=(X(I+1)+X(I+2))/2.0	PH1 1990
3360 CONTINUE	PH1 2000
3361 IF(VEL)9911,10000,3363	PH1 2010
3363 VEL=0.0	PH1 2020
GO TO 3301	PH1 2030
C ERROR	PH1 2040
9902 NK=3305	PH1 2050
GO TO 9999	PH1 2060
9903 NK=3306	PH1 2070
GO TO 9999	PH1 2080
9904 NK=3310	PH1 2090
GO TO 9999	PH1 2100
9905 NK=3316	PH1 2110
GO TO 9999	PH1 2120
9906 NK=3320	PH1 2130
GO TO 9999	PH1 2140
9907 NK=3328	PH1 2150
GO TO 9999	PH1 2160
9911 NK=3361	PH1 2170
9999 NR=3	PH1 2180
CALL DUMP	PH1 2190
C IF SN(NOT=0.) ANY NEGATIVE ENERGIES WILL	
C REMAIN. IF=0, CODE WILL TRY ANOTHER PASS	
C WITH A SMALLER DT.	
10000 IF(SN)7030,7031,7030	PH1 2200
7031 IF(UT)7020,7030,7010	PH1 2210
C NEGATIVE ENERGIES HAVE OCCURED, INTEGRATE	
C BACK TO CYCLE N WITH (-DT).	
7010 UT=-1.0	PH1 2220
DT=-DT	PH1 2230
C YOU NOW HAVE INTEGRATED BACK TO CYCLE N. NOW	
C INTEGRATE TO CYCLE N+1 WITH NEW DT(STORED IN UU).	
GO TO 8000	PH1 2240
7020 UT=0.0	PH1 2250
DT=UU	PH1 2260
NR=DT/TRAD+1.0	PH1 2270
WS=NR	PH1 2280

116.

PH1 2290
PH1 2300
PH1 2310

PH1 2320
PH1 2330
PH1 2340
PH1 2350
PH1 2360
PH1 2370
PH1 2380
PH1 2390

TRAD=DT/WS

DTNA=DT

GO TO 8000

C INCREASE ACTIVE GRID COUNTERS IF NEEDED.

7030 I1=I1+NRC

I2=I2+NRT

IF(I1-IMAX)6100,6100,6200

6200 I1=IMAX

6100 IF(I2-JMAX)6201,6201,6202

6202 I2=JMAX

6201 RETURN

END

\$IBFTC PH2 LIST,DECK,REF

117.

SUBROUTINE PH2

PH2 0010

C Z(110)= CRITICAL ENERGY(BETWEEN GAS AND CONDENSED STATE)
C Z(111)= INITIAL DENSITY
C Z(112)= INITIAL VELOCITY OF PELLET
C Z(113)= EPSILONICS FOR EMPTYING PELLET ON BASIS OF VELOCITY
C TOZONE = MINIMUM DENSITY FOR MASS FLOW

PH2 0900

PH2 0980

C
C
C AMPY=MASS ACROSS TOP BOUNDARY OF CELL
C AMUT=RADIAL MOMENTA OF THIS MASS
C AMVT=AXIAL MOMENTA OF THIS MASS
C DELET=TOTAL SPECIFIC ENERGY OF THIS MASS
C AMMP=MASS ACROSS RIGHT BOUNDARY OF CELL
C AMUR=RADIAL MOMENTA OF THIS MASS
C AMVR=AXIAL MOMENTA OF THIS MASS
C DELER=TOTAL SPECIFIC ENERGY OF THIS MASS
C AMMY=MASS ACROSS BOTTOM BOUNDARY OF CELL
C AMMU=RADIAL MOMENTA OF THIS MASS
C AMMV=AXIAL MOMENTA OF THIS MASS
C DELEB=TOTAL SPECIFIC ENERGY OF THIS MASS
C GAMC=MASS ACROSS LEFT BOUNDARY OF CELL
C FLEFT=RADIAL MOMENTA OF THIS MASS
C YAMC=AXIAL MOMENTA OF THIS MASS
C SIGC=TOTAL SPECIFIC ENERGY OF THIS MASS
C =====

PH2 0990

PH2 1010

NRT=0

PH2 1020

NRC=0

PH2 1030

REZ=0.0

PH2 1040

CALL SLITE (0)

PH2 1050

PIDTS=1.0/(PIDY*DT)

PH2 1060

101 DO 103 J=1,JMAX

PH2 1070

102 GAMC(J)=0.0

PH2 1080

FLEFT(J)=0.0

PH2 1090

YAMC(J)=0.0

PH2 1100

SIGC(J)=0.0

PH2 1110

103 CONTINUE

PH2 1120

104 DO 547 I=1,I1

PH2 1130

J=I

PH2 1140

105 K=I+1

PH2 1150

80 IF(AMX(K))9900,97,81

PH2 1160

81 IF(V(K))82,97,97

PH2 1170

97 AMMV=0.0

PH2 1180

GO TO 98

PH2 1190

82 AMMY=AMX(K)*V(K)*DT/DY(J)

PH2 1200

83 IF(AMMY+AMX(K))84,85,85

PH2 1210

84 AMMY=-AMX(K)

PH2 1220

85 IF(CVIS))06,99,99

PH2 1230

C BOTTOM BOUNDARY IS TRANSMITTIVE, MATERIAL IS MOVING

..8.

C	OUT, REMOVE ITS ENERGY FROM ETH.	
106	AMMU=AMMY*U(K)	PH2 1240
	AMMV=AMMY*V(K)	PH2 1250
	DELEB=AIX(K)+(U(K)**2+V(K)**2)/2.0	PH2 1260
	WS=(U(K)**2+V(K)**2)/2.0	PH2 1270
	ETH=ETH+AMMY*(AEX(K)+WS)	PH2 1280
	GO TO 107	PH2 1290
C	BOTTOM BOUNDARY IS REFLECTIVE, NET MOMENTA CHANGE	
C	IN Z DIRECTION IS 2 MV.	
99	AMMV=2.0*AMMY*V(K)	PH2 1300
98	AMMY=0.0	PH2 1310
	AMMU=0.0	PH2 1320
	DELEB=0.0	PH2 1330
C	BEGIN DO LOOP IN J(Z) DIRECTION.	
107	DO 546 J=1,12	PH2 1340
108	L=K+IMAX	PH2 1350
	I=I	PH2 1360
	J=J	PH2 1370
	AREA=0.0	PH2 1380
	VEL=0.0	PH2 1390
	FS=0.0	PH2 1400
210	IF(JMAX-J)211,211,207	PH2 1410
211	VEL=1.0	PH2 1420
	GO TO 208	PH2 1430
207	IF(AMX(L))215,215,214	PH2 1440
214	IF(AMX(K))216,216,209	PH2 1450
216	VABOVE=V(L)	PH2 1460
	GO TO 212	PH2 1470
215	IF(AMX(K))205,205,208	PH2 1480
205	VABOVE=0.0	PH2 1490
	GO TO 212	PH2 1500
208	VABOVE=V(K)	PH2 1510
	GO TO 212	PH2 1520
209	VABOVE=(V(K)+V(L))/2.0	PH2 1530
212	CONTINUE	PH2 1540
	I=I	PH2 1550
	J=J	PH2 1560
	FS=0.0	PH2 1570
404	IF(IMAX-I)412,412,405	PH2 1580
405	IF(AMX(K+1))411,411,409	PH2 1590
409	IF(AMX(K))410,410,407	PH2 1600
410	URR=U(K+1)	PH2 1610
	GO TO 408	PH2 1620
411	IF(AMX(K))403,403,406	PH2 1630
403	URR=0.0	PH2 1640
	GO TO 408	PH2 1650
C	WE ARE AT THE RIGHT BOUNDARY OF THE GRID, THE	
C	BOUNDARY CONDITION HERE IS TRANSMITTIVE.	
412	FS=1.0	PH2 1660
406	URR=U(K)	PH2 1670

GO TO 408	PH2 1680
407 URR=(U(K)+U(K+1))/2.0	PH2 1690
408 CONTINUE	PH2 1700
109 IF(AREA)9901,301,547	PH2 1710
301 IF(VABOVE)300,304,302	PH2 1720
302 IF(AMX(K))9900,304,8800	PH2 1730
8800 IF(J-1)9900,303,8801	PH2 1740
8801 KP=K-IMAX	PH2 1750
IF(AMX(KP))9900,8803,303	PH2 1760
C A CHECK HERE TO INSURE THAT THE BOTTOM ZONES	
C OF THE PROJECTILE EMPTY (FOR HYPERVELOCITY) UP UNTIL	
C THE INITIAL VELOCITY CHANGES DUE TO THE SHOCK.	
8803 IF(ABS(VABOVE-Z(112))/Z(112)-Z(113))306,303,303	PH2 1770
303 M=K	PH2 1780
JJ=J	PH2 1790
GO TO 307	PH2 1800
304 AMPY=0.0	PH2 1810
308 AMUT=0.0	PH2 1820
AMVT=0.0	PH2 1830
DELET=0.0	PH2 1840
GO TO 501	PH2 1850
300 IF(VEL)9901,305,304	PH2 1860
305 IF(AMX(L))9903,304,306	PH2 1870
306 M=L	PH2 1880
JJ=J+1	PH2 1890
307 IF(VEL)6130,6130,6140	PH2 1900
6130 WSA=(V(K)+V(L))/2.0	PH2 1910
WSB=1.0+(V(L)-V(K))/(DY(JJ)*SBOUND)*DT	PH2 1920
VABOVE=WSA/WSB	PH2 1930
C CALCULATE THE MASS FLUX AT THE TO. OF CELL K.	
6140 AMPY=AMX(M)*VABOVE/DY(JJ)*DT	PH2 1940
501 IF(URR)500,504,502	PH2 1950
502 IF(AMX(K))9900,504,503	PH2 1960
503 M=K	PH2 1970
N=I	PH2 1980
GO TO 508	PH2 1990
504 AMMP=0.0	PH2 2000
AMUR=0.0	PH2 2010
AMVR=0.0	PH2 2020
DELER=0.0	PH2 2030
GO TO 1	PH2 2040
500 IF(FS)9905,506,504	PH2 2050
506 IF(AMX(K+1))9904,504,507	PH2 2060
507 M=K+1	PH2 2070
N=I+1	PH2 2080
508 IF(FS)6100,6100,6110	PH2 2090
6100 WSA=(U(K)+U(K+1))/2.0	PH2 2100
WSB=1.0+(U(K+1)-U(K))/(DX(N)*SBOUND)*DT	PH2 2110
URR=WSA/WSB	PH2 2120
C CALCULATE THE MASS FLUX AT THE RIGHT OF CELL K.	

6110 DEN=AMX(M)/TAU(N)	PH2 2130
AMMP=DEN/PIDTS*X(1)/.5*URR	
1 IF(AMMP)16,74,8820	PH2 2150
8820 IF(GAMC(J))74,74,8821	PH2 2160
8821 IF(FS)6120,6120,74	PH2 2170
6120 IF(AMX(K+1))9903,8822,74	PH2 2180
8822 IF(AMX(K)/(TAU(I)*DY(J))-Z(111))8823,74,74	PH2 2190
8823 IF(AIX(K)-Z(110))8824,74,74	PH2 2200
8824 WS=GAMC(J)+AMX(K)-TAU(I)*DY(J)*Z(111)	PH2 2210
IF(WS)8826,8826,8825	PH2 2220
8825 AMMP=WS	PH2 2230
GO TO 74	PH2 2240
8826 AMMP=0.0	PH2 2250
74 JTAG=0	PH2 2260
C BEGIN CHECKING TO SEE IF THEIR IS ANY	
C PREFERENTIAL MASS FLUX BECAUSE OF CHOICE OF	
C INDEXING DIRECTION.	
2 IF(AMPY)3,4,4	PH2 2270
C TOP FLUX IS INTO CELL K.	
3 ITAG=1	PH2 2280
WSB=AMPY	PH2 2290
AMPY=0.0	PH2 2300
GO TO 64	PH2 2310
4 ITAG=0	PH2 2320
64 IF(AMMY)9,5,5	PH2 2330
C BOTTOM FLUX IS INTO CELL K.	
5 IF(GAMC(J))7,6,6	PH2 2340
C LEFT FLUX IS INTO CELLK.	
6 WS=AMX(K)	PH2 2350
GO TO 11	PH2 2360
C LEFT FLUX IS OUT.	
7 WS=AMX(K)+GAMC(J)	PH2 2370
GO TO 11	PH2 2380
C BOTTOM FLUX IS OUT OF CELL K.	
9 IF(GAMC(J))10,8,8	PH2 2390
C LEFT FLUX IS INTO CELL K.	
8 WS=AMX(K)+AMMY	PH2 2400
GO TO 11	PH2 2410
C LEFT FLUX IS OUT OF CELL K.	
10 WS=AMX(K)+GAMC(J)+AMMY	PH2 2420
11 WSA=AMPY+AMMP	PH2 2430
12 IF(WSA-WS)75,75,13	PH2 2440
C CHANGE TOP AND RIGHT FLUX TO BE THE	
C OLD FLUX TIMES THE MASS OF THE CELL/THE SUM	
C OF THE OLD FLUXES.	
13 AMPY=AMPY*WS/WSA	PH2 2450
AMMP=AMMP*WS/WSA	PH2 2460
75 IF(JTAG)14,73,14	PH2 2470
73 WSC=AMMP	PH2 2480
14 IF(ITAG)15,7000,15	PH2 2490

15	AMPY=WSB	PH2 2500
	ITAG=0	PH2 2510
C	GO CHECK CELL ABOVE.	
	GO TO 40	PH2 2520
C	RIGHT FLUX IS INTO CELL K.	
16	IF(FS) 76, 17, 76	PH2 2530
76	WSC=AMMP	PH2 2540
C	I=IMAX, SO CHECK CELL ABOVE K.	
	GO TO 40	PH2 2550
17	IF(I+1-IMAX) 19, 19, 9908	PH2 2560
18	URRR=U(K+1)/2.0	PH2 2570
	GO TO 20	PH2 2580
19	URRR=(U(K+1)+U(K+2))/2.0	PH2 2590
20	IF(URRR) 39, 39, 21	PH2 2600
C	FLUX IS OUT OF THE RIGHT OF CELL(K+1).	
21	URRR=URRR/TAU(I+1)*AMX(K+1)/PID15*X(I+1)/.5	
22	IF(J-1) 9909, 23, 24	PH2 2620
23	VBLO=V(K+1)/2.0	PH2 2630
	GO TO 26	PH2 2640
24	KP=K+1-IMAX	PH2 2650
	VBLO=(V(K+1)+V(KP))/2.0	PH2 2660
26	IF(VBLO) 25, 38, 38	PH2 2670
C	FLUX IS OUT OF THE BOTTOM OF CELL(K+1).	
25	VBLO=AMX(K+1)/DY(J)*VBLO*DT	PH2 2680
27	IF(VEL) 28, 29, 28	PH2 2690
28	VAB=V(K+1)/2.0	PH2 2700
	GO TO 31	PH2 2710
29	KP=K+IMAX+1	PH2 2720
	VAB=(V(K+1)+V(KP))/2.0	PH2 2730
31	IF(VAB) 36, 36, 30	PH2 2740
C	FLUX IS OUT OF TOP.	
30	VAB=AMX(K+1)/DY(J)*VAB*DT	PH2 2750
32	WS=AMX(K+1)	PH2 2760
33	WSA=URRR-AMMP-VBLO+VAB	PH2 2770
34	IF(WSA-WS) 77, 77, 35	PH2 2780
35	AMMP=AMMP*WS/WSA	PH2 2790
77	JTAG=1	PH2 2800
	WSC=AMMP	PH2 2810
	AMMP=0.0	PH2 2820
	GO TO 2	PH2 2830
C	FLUX AT TOP IS INTO CELL (K+1).	
36	WS=AMX(K+1)	PH2 2840
37	WSA=URRR-AMMP-VBLO	PH2 2850
	GO TO 34	PH2 2860
C	FLUX IS IN FROM BOTTOM INTO CELL (K+1).	
38	VBLO=0.0	PH2 2870
	GO TO 27	PH2 2880
C	FLUX IS INTO CELL (K+1) FROM RIGHT.	
39	URRR=C.0	PH2 2890
	GO TO 22	PH2 2900

C	RIGHT FLUX OUT OF CELL (K) IS POSITIVE AND TOP	
C	FLUX IS COMING INTO CELL (K) FROM (K+IMAX).	
40	IF(VEL)7000,41,7000	PH2 2910
41	IF(FS)42,43,42	PH2 2920
C	WE ARE AT THE RIGHT BOUNDARY OF THE GRID.	
42	KP=K+IMAX	PH2 2930
	URT=U(KP)/2.0	PH2 2940
	GO TO 45	PH2 2950
43	KP=K+IMAX	PH2 2960
	URT=(U(KP)+U(KP+1))/2.0	PH2 2970
45	IF(URT)46,46,70	PH2 2980
C	FLUX AT RIGHT (CELL M) IS NEGATIVE.	
46	URT=0.0	PH2 2990
	GO TO 47	PH2 3000
70	KP=K+IMAX	PH2 3010
	URT=URT/TAU(I)*AMX(KP)/PIDTS*X(I)/.5	
C	FLUX AT RIGHT (CELL M) IS POSITIVE.	
47	IF(J+1-JMAX)48,49,9910	PH2 3030
48	KP=K+IMAX	PH2 3040
	KL=KP+IMAX	PH2 3050
	VABT=(V(KP)+V(KL))/2.0	PH2 3060
	GO TO 51	PH2 3070
49	KP=K+IMAX	PH2 3080
	KL=KP+IMAX	PH2 3090
	VABT=V(KP)/2.0	PH2 3100
51	IF(VABT)8810,71,72	PH2 3110
C	FLUX IS IN FROM TOP OF CELL M.	
8810	IF(AMX(K))9903,8811,71	PH2 3120
C	CHECK FOR SOLID OR VAPOR.	
8811	IF(AMX(KP)/(TAU(I)*DY(J+1))-Z(111))8812,71,71	PH2 3130
8812	IF(AIX(KP)-Z(110))8813,71,71	PH2 3140
8813	VABT=VABT*AMX(KL)/DY(J+2)*DT	PH2 3150
8814	WS=-VABT+AMX(KP)-TAU(I)*DY(J+1)*Z'(111)	PH2 3160
8815	IF(WS)8817,8817,8816	PH2 3170
8816	AMPY=-WS	PH2 3180
	GO TO 71	PH2 3190
8817	AMPY=0.0	PH2 3200
71	VABT=0.0	PH2 3210
	GO TO 60	PH2 3220
72	VABT=VABT*AMX(KP)/DY(J+1)*DT	PH2 3230
52	IF(GAMC(J+1))54,53,53	PH2 3240
53	WS=AMX(KP)	PH2 3250
	GO TO 55	PH2 3260
54	WS=AMX(KP)+GAMC(J+1)	PH2 3270
55	WSA=VABT-AMPY+URT	PH2 3280
	GO TO 57	PH2 3290
60	IF(GAMC(J+1))6,61,59	PH2 3300
61	WS=AMX(KP)+GAMC(J+1)	PH2 3310
	GO TO 58	PH2 3320
59	WS=AMX(KP)	PH2 3330

58	WSA=-AMPY+URT	PH2 3340
57	IF(WSA-WS)7000,7000,56	PH2 3350
56	AMPY=AMPY*WS/WSA	PH2 3360
	GO TO 7000	PH2 3370
7000	AMMP=WSC	PH2 3380
309	IF(AMPY)8834,8831,8833	PH2 3390
8833	IF(JMAX-J)9911,318,8835	PH2 3400
8835	KP=K+IMAX	PH2 3410
8836	IF(AMX(KP))9900,8837,318	PH2 3420
C	**** NOTE ****	
C	ACROSS FREE SURFACE, HOLD UP MASS FLUX	
C	UNLESS THIS MASS PRODUCES A DENSITY GREATER THAN TOZONE.	
C	*****	
8837	IF(AMPY/(TAU(I)*DY(J))-TOZONE)8838,318,318	PH2 3430
8838	AMPY=0.0	PH2 3440
	GO TO 8831	PH2 3450
8834	IF(J-1)9911,325,8839	PH2 3460
8839	IF(AMX(K))9900,8840,325	PH2 3470
8840	IF(-AMPY/(TAU(I)*DY(J))-TOZONE)8841,325,325	
8841	AMPY=0.0	PH2 3490
	GO TO 8831	PH2 3500
318	DELM=GAMC(J)+AMMY-AMPY	PH2 3510
322	IF(VEL)9901,324,323	PH2 3520
323	WS=U(K)**2+V(K)**2	PH2 3530
C	MATERIAL HAS LEFT THE TOP, TRIGGER REZONE	
C	FLAG, REMOVE ITS ENERGY FROM ETH(TOTAL ENERGY OF SYSTEM).	
	ETH=ETH-AMPY*(AIX(K)+WS/2.0)	PH2 3540
	IF(AMPY/(TAU(I)*DY(J))-TOZONE)324,324,6900	PH2 3550
6900	REZ=1.0	PH2 3560
324	AMUT=AMPY*U(K)	PH2 3570
	AMVT=AMPY*V(K)	PH2 3580
	GO TO 326	PH2 3590
325	CONTINUE	PH2 3600
8831	AMUT=AMPY*U(L)	PH2 3610
	AMVT=AMPY*V(L)	PH2 3620
	DELM=GAMC(J)-AMPY+AMMY	PH2 3630
326	IF(AMPY)327,328,328	PH2 3640
327	DELET=AIX(L)+(U(L)**2+V(L)**2)/2.0	PH2 3650
	GO TO 333	PH2 3660
328	IF(AMMY)329,330,330	PH2 3670
329	DELET=DELEB	PH2 3680
	GO TO 333	PH2 3690
330	IF(GAMC(J))331,332,332	PH2 3700
331	DELET=SIGC(J)	PH2 3710
	GO TO 333	PH2 3720
332	DELET=AIX(K)+(U(K)**2+V(K)**2)/2.0	PH2 3730
C	SUM UP RADIAL MOMENTA FOR ALL FLUXES EXCEPT	
C	THE RIGHT AND STORE IN SIGMU.	
333	SIGMU=FLEFT(J)+AMMU-AMUT	PH2 3740
C	SUM UP AXIAL MOMENTA FOR ALL FLUXES EXCEPT THE	

124.

C	RIGHT AND STORE IN SIGMV.	
	SIGMV=YAMC(J)+AMMV-AMVT	PH2 3750
C	SUM UP TOTAL ENERGY CARRIED BY THESE FLUXES	
C	EXCEPT THE RIGHT FLUX AND STORE IN DELEK.	
	DELEK=GAMC(J)*SIGC(J)+AMMY*DELES-AMPY*DELET	PH2 3760
509	IF(AMMP)8843,518,8844	PH2 3770
8844	IF(IMAX-I)9911,518,8845	PH2 3780
8845	IF(AMX(K+1))9900,8846,518	PH2 3790
8846	IF(AMMP/(TAU(I)*DY(J))-TOZONE)8847,518,518	PH2 3800
8847	AMMP=0.0	PH2 3810
	GO TO 518	PH2 3820
8843	IF(I-1)9911,512,8848	PH2 3830
8848	IF(AMX(K))9900,8849,512	PH2 3840
8849	IF(-AMMP/(TAU(I)*DY(J))-TOZONE)8850,512,512	
8850	AMMP=0.0	PH2 3860
	GO TO 518	PH2 3870
512	DELM=DELM-AMMP+AMX(K)	PH2 3880
513	CONTINUE	PH2 3890
514	CONTINUE	PH2 3900
8828	AMUR=AMMP*U(K+1)	PH2 3910
	AMVR=AMMP*V(K+1)	PH2 3920
	GO TO 525	PH2 3930
518	DELM=DELM-AMMP+AMX(K)	PH2 3940
521	CONTINUE	PH2 3950
522	IF(FS)9905,524,523	PH2 3960
523	WS=U(K)**2+V(K)**2	PH2 3970
	ETH=ETH-AMMP*(AIX(K)+WS/2.0)	PH2 3980
	IF(AMMP/(TAU(I)*DY(J))-TOZONE)524,524,6901	PH2 3990
6901	REZ=1.0	PH2 4000
524	AMUR=AMMP*U(K)	PH2 4010
	AMVR=AMMP*V(K)	PH2 4020
525	SIGMU=SIGMU-AMUR	PH2 4030
	SIGMV=SIGMV-AMVR	PH2 4040
526	TIC=0.0	PH2 4050
527	IF(AMMP)528,529,529	PH2 4060
528	DELER=AIX(K+1)+(U(K+1)**2+V(K+1)**2)/2.0	PH2 4070
	GO TO 537	PH2 4080
529	IF(AMMY)530,531,531	PH2 4090
530	DELER=DELEB	PH2 4100
	GO TO 536	PH2 4110
531	IF(GAMC(J))532,533,533	PH2 4120
532	DELER=SIGC(J)	PH2 4130
	GO TO 536	PH2 4140
533	IF(AMPY)535,535,534	PH2 4150
534	DELER=DELET	PH2 4160
	GO TO 536	PH2 4170
535	DELER=AIX(K)+(U(K)**2+V(K)**2)/2.0	PH2 4180
536	TIC=1.0	PH2 4190
537	DELEK=DELEK-AMMP*DELER	PH2 4200
538	IF(TIC)9907,539,550	PH2 4210

550 WS=DELER	PH2 4220
GO TO 999	PH2 4230
539 WS=AIX(K)+(U(K)**2+V(K)**2)/2.0	PH2 4240
999 IF(DELM)998,543,540	PH2 4250
998 IF(AMX(K)*1.E-6+DELM)9906,997,997	PH2 4260
997 DELM=0.0	PH2 4270
GO TO 543	PH2 4280
C ENK=TOTAL ENERGY OF CELL (K) + ENERGY THAT	
C HAS BEEN ADDED AND LOST.	
540 ENK=AMX(K)*WS+DELEK	PH2 4290
C BY CONSERVING AXIAL MOMENTA, CALCULATE THE NEW	
C AXIAL VELOCITY COMPONENT FOR CELL K.	
541 U(K)=(SIGMU+AMX(K)*U(K))/DELM	PH2 4300
C BY CONSERVING RADIAL MOMENTA, CALCULATE THE NEW	
C RADIAL VELOCITY COMPONENT FOR CELL K.	
601 V(K)=(SIGAV+AMX(K)*V(K))/DELM	PH2 4310
IF(I-I1)603,6604,6604	PH2 4320
6604 IF(U(K))6605,6606,6605	PH2 4330
6605 NRC=1	PH2 4340
6606 IF(V(K))6607,6608,6607	PH2 4350
6607 NRC=1	PH2 4360
6608 IF(AIX(K))6609,6610,6609	PH2 4370
6609 NRC=1	PH2 4380
6610 CONTINUE	PH2 4390
603 WS=U(K)**2+V(K)**2	PH2 4400
C BY CONSERVING BOTH TOTAL ENERGY AND	
C MOMENTA, CALCULATE THE NEW SPECIFIC	
C INTERNAL ENERGY FOR CELL K.	
542 AIX(K)=ENK/DELM-WS/2.0	PH2 4410
543 AMX(K)=DELM	PH2 4420
IF(AMX(K))9900,2007,544	PH2 4430
2007 AIX(K)=0.0	PH2 4440
U(K)=0.0	PH2 4450
V(K)=0.0	PH2 4460
P(K)=0.0	PH2 4470
C THE RIGHT VALUES OF CELL (K) BECOME THE LEFT	
C VALUES OF CELL (K+1).	
544 GAMC(J)=AMMP	PH2 4480
FLEFT(J)=AMUR	PH2 4490
YAMC(J)=AMVR	PH2 4500
SIGC(J)=DELER	PH2 4510
C THE TOP VALUES OF CELL(K) BECOME THE	
C BOTTOM VALUES FOR CELL (K+IMAX).	
545 AMMY=AMPY	PH2 4520
AMMU=AMUT	PH2 4530
AMMV=AMVT	PH2 4540
DELEB=DELET	PH2 4550
546 K=K+IMAX	PH2 4560
LL=K-IMAX	PH2 4570
IF(U(LL))6500,6600,6500	PH2 4580


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6500 NRT=1
6600 IF(V(LL))6601,6602,6601
6601 NRT=1
6602 IF(AIX(L))6611,547,6611
6611 NRT=1
  547 CONTINUE
      I1=I1+NRC
      I2=I2+NRT
      IF(IMAX-I1)6700,6701,6702
6700 I1=IMAX
6701 CONTINUE
6702 IF(JMAX-I2)6800,6801,6802
6800 I2=JMAX
6801 CONTINUE
6802 GO TO 548
9901 NK=300
      GO TO 9999
9900 NK=302
      GO TO 9999
9903 NK=305
      GO TO 9999
9904 NK=506
      GO TO 9999
9905 NK=500
      GO TO 9999
9906 NK=513
      GO TO 9999
9911 NK=8833
      GO TO 9999
9908 NK= 17
      GO TO 9999
9909 NK= 22
      GO TO 9999
9910 NK= 47
      GO TO 9999
9907 NK=538
9999 NR=4
      CALL DUMP
  548 SUM=0.0
2005 DO 2001 I=1,I1
      K=I+1
      DO 2000 J=1,I2
        IF(AMX(K))2000,2000,2009
C      IF ANY RHO (CELL DENSITY) IS LESS THAN TOZONE,
C      SET THE MASS TO ZERO, AND TALLY THE
C      MOMENTAS AND ENERGIES IN THE Z STORAGE, ALSO
C      CHECK FOR NEGATIVE INTERNAL ENERGIES, IF
C      WE FIND SOME, SET THEM TO ZERO AFTER
C      SUBTRACTING THEM FROM ETH..
2009 IF(AMX(K)/(TAU(I)*DY(J))-TOZONE)2010,2008,2008

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PH2 4590
PH2 4600
PH2 4610
PH2 4620
PH2 4630
PH2 4640
PH2 4650
PH2 4660
PH2 4670
PH2 4680
PH2 4690
PH2 4700
PH2 4710
PH2 4720
PH2 4730
PH2 4740
PH2 4750
PH2 4760
PH2 4770
PH2 4780
PH2 4790
PH2 4800
PH2 4810
PH2 4820
PH2 4830
PH2 4840
PH2 4850
PH2 4860
PH2 4870
PH2 4880
PH2 4890
PH2 4900
PH2 4910
PH2 4920
PH2 4930
PH2 4940
PH2 4950
PH2 4960
PH2 4970
PH2 4980
PH2 4990
PH2 5000
PH2 5010

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PH2 5020
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2010 WS=(U(K)**2+V(K)**2)/2.0
      Z(100)=Z(100)+AMX(K)
      WS=AMX(K)*(AIX(K)+WS)
      Z(101)=Z(101)+WS
      ETH=ETH-WS
      Z(102)=Z(102)+AMX(K)*U(K)
      Z(103)=Z(103)+AMX(K)*V(K)
      AMX(K)=0.0
      AIX(K)=0.0
      P(K)=0.0
      U(K)=0.0
      V(K)=0.0
      GO TO 2000
2008 IF(AIX(K))2004,2000,2000
2004 SUM=SUM+AIX(K)*AMX(K)
      AIX(K)=0.0
2000 K=K+IMAX
2001 CONTINUE
2003 ETH=ETH-SUM
      Z(104)=Z(104)+SUM
8000 IF(REZ)8001,8001,8002
8002 IF(REZFCT)8004,8004,8003
8004 REZ=0.
      GO TO 8001
8003 CALL REZONE
8001 RETURN
      END

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PH2 5030
PH2 5040
PH2 5050
PH2 5060
PH2 5070
PH2 5080
PH2 5090
PH2 5100
PH2 5110
PH2 5120
PH2 5130
PH2 5140
PH2 5150
PH2 5160
PH2 5170
PH2 5180
PH2 5190
PH2 5200
PH2 5210
PH2 5220

```

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PH2 5260
PH2 5270

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\$IBFTC ES	LIST,DECK,REF	
	SUBROUTINE ES	ES 0010
C		ES 0760
C		ES 0900
C	METALLIC EQUATION OF STATE, SEE	
C	GA-3216 REPORT.	
C		ES 0980
10	RHOW=AMX(K)/(TAU(I)*DY(J))	ES 0990
	ETA=RHOW/Z(115)	ES 1000
	VOW=1.0/ETA	ES 1010
11	P1=AIX(K)*RHOW*Z(116)	ES 1020
12	P2=(Z(115)*TAU(I)*DY(J))**2*AIX(K)	ES 1030
13	P3=AMX(K)**2*Z(117)	ES 1040
14	P4=Z(118)/(P2/P3+1.0)*AIX(K)*RHOW	ES 1050
15	P5=Z(119)*(ETA-1.0)	ES 1060
16	IF(ETA-1.0)50,100,100	ES 1070
50	IF(VOW-Z(120))55,55,75	ES 1080
55	IF(AIX(K)-Z(122))100,100,75	ES 1090
75	P7=Z(123)*(VOW-1.0)	ES 1100
	IF(P7-88.0)4002,4002,4003	ES 1110
4003	P7=88.0	ES 1120
4002	CONTINUE	ES 1130
	P8=EXP(P7)	ES 1140
	P9=1.0/P8	ES 1150
	P10=Z(124)*((VOW-1.0)**2)	ES 1160
	IF(P10-88.0)4000,4000,4001	ES 1170
4001	P10=88.0	ES 1180
4000	CONTINUE	ES 1190
	P11=EXP(P10)	ES 1200
	P12=1.0/P11	ES 1210
	P(K)=P1+(P4+P5*P9)*P12	ES 1220
	GO TO 119	
100	P6=Z(126)*((ETA-1.0)**2)	ES 1230
	P(K)=P1+P4+P5+P6	ES 1240
119	IF(P(K))999,999,200	ES 1250
200	WSGX=.5	ES 1260
	GO TO 500	ES 1270
999	P(K)=0.0	ES 1280
	WSGX=.5+Z(125)	ES 1290
	GO TO 500	ES 1300
500	RETURN	ES 1310
	END	ES 1320

\$IBFTC ES	LIST,DECK,REF	
	SUBROUTINE ES	ES 0010
C	** POLYTROPIC EQUATION OF STATE **	ES 0980
	P(K)=AMX(K)*AIX(K)/GAMX/(DY(J)*TAU(I))	ES 0990
	WSGX=GMAX-1.0	ES 1000
	RETURN	ES 1010
	END	ES 1020

3IBFTC REZONE LIST,DECK,REF
SUBROUTINE REZONE

REZ00010
REZ00980

C
C CONSERVE MOMENTUM AND TOTAL ENERGY, INCREASE
C ALL LINEAR DIMENSIONS BY 2. (THUS 4 CELLS
C IN THE OLD GRID ARE COMBINED INTO 1 FOR
C THE NEW GRID.)
NIMAX=IMAX/2
NJMAX=JMAX/2
DO 10 J=1,NJMAX
K=(J-1)*NIMAX+2
L=(J-1)*2*IMAX+2
DO 11 I=1,NIMAX
M=L+IMAX
12 WSA=AMX(L)+AMX(M)+AMX(L+1)+AMX(M+1)
WSB=AMX(L)*(U(L)**2+V(L)**2)+AMX(M)*(U(M)
1**2+V(M)**2)+AMX(L+1)*(U(L+1)**2+V(L+1)**2)
2+AMX(M+1)*(U(M+1)**2+V(M+1)**2)
U(K)=(U(L)*AMX(L)+U(M)*AMX(M)+U(L+1)*AMX(L+1)+
1U(M+1)*AMX(M+1))/WSA
V(K)=(V(L)*AMX(L)+V(M)*AMX(M)+V(L+1)*AMX(L+1)+
1V(M+1)*AMX(M+1))/WSA
AIX(K)=AIX(L)*AMX(L)+AIX(M)*AMX(M)+AIX(L+1)*
1AMX(L+1)+AMX(M+1)*AIX(M+1)
AMX(K)=WSA
WS=U(K)**2+V(K)**2
E=AIX(K)+WSB/2.0
AIX(K)=E/AMX(K)-.5*WS
IF(K-2)14,14,13
C SET CELL QUANTITIES OF OLD GRID TO ZERO.
13 AMX(L)=0.0
U(L)=0.0
V(L)=0.0
AIX(L)=0.0
P(L)=0.0
AMX(M)=0.0
U(M)=0.0
V(M)=0.0
AIX(M)=0.0
P(M)=0.0
AMX(L+1)=0.0
U(L+1)=0.0
V(L+1)=0.0
AIX(L+1)=0.0
P(L+1)=0.0
AMX(M+1)=0.0
U(M+1)=0.0
V(M+1)=0.0
AIX(M+1)=0.0
P(M+1)=0.0

REZ00990
REZ01000
REZ01010
REZ01020
REZ01030
REZ01040
REZ01050
REZ01060
REZ01070
REZ01080
REZ01090
REZ01100
REZ01110
REZ01120
REZ01130
REZ01140
REZ01150
REZ01160
REZ01170
REZ01180
REZ01190
REZ01200

REZ01210
REZ01220
REZ01230
REZ01240
REZ01250
REZ01260
REZ01270
REZ01280
REZ01290
REZ01300
REZ01310
REZ01320
REZ01330
REZ01340
REZ01350
REZ01360
REZ01370
REZ01380
REZ01390
REZ01400

199.

14	K=K+1	REZ01410
	L=L+2	REZ01420
11	CONTINUE	REZ01430
10	CONTINUE	REZ01440
C	CALCULATE NEW DY AND Y (JMAX OF THEM).	
18	DO 999 J=1,JMAX	REZ01450
	DY(J)=DY(J)*2.0	REZ01460
999	CONTINUE	REZ01470
	DO 99 J=1,JMAX	REZ01480
	Y(J)=Y(J-1)+DY(J)	REZ01490
99	CONTINUE	REZ01500
16	DX(1)=2.0*DX(1)	REZ01510
	X(1)=DX(1)	REZ01520
	WS=X(1)**2	REZ01530
	TAU(1)=PIDY*WS	REZ01540
C	CALCULATE NEW DX AND X, AND TAU(IMAX OF THEM)	
17	DO 98 I=2,IMAX	REZ01550
	X(I)=X(I-1)+DX(I)	REZ01560
	DX(I)=DX(I)	REZ01570
	WSA=X(I)**2	REZ01580
	TAU(I)=PIDY*(WSA-WS)	REZ01590
	WS=WSA	REZ01600
98	CONTINUE	REZ01610
	IMAX=NIMAX	REZ01620
	JMAX=NJMAX	REZ01630
C	PREPARE NOW TO SHUFFLE THE X ARRAYS SUCH	
C	AS TO PRESERVE K=(J-1)*IMAX+1+1.	
	DO 20 N=1,JMAX	REZ01640
	J=JMAX+1-N	REZ01650
	K=(J-1)*IMAX+1+IMAX	REZ01660
	L=(J-1)*(IMAX+IMAX)+1+IMAX	REZ01670
	DO 21 I=1,IMAX	REZ01680
1000	AMX(L)=AMX(K)	REZ01690
	AIX(L)=AIX(K)	REZ01700
	U(L)=U(K)	REZ01710
	V(L)=V(K)	REZ01720
	P(L)=P(K)	REZ01730
	IF(J-1)1002,1002,1001	REZ01740
1001	AMX(K)=0.0	REZ01750
	AIX(K)=0.0	REZ01760
	V(K)=0.0	REZ01770
	U(K)=0.0	REZ01780
	P(K)=0.0	REZ01790
1002	K=K-1	REZ01800
	L=L-1	REZ01810
21	CONTINUE	REZ01820
20	CONTINUE	REZ01830
	IMAX=NIMAX*2	REZ01840
	JMAX=NJMAX*2	REZ01850
	II=NIMAX+1	

```

      JJ=KJMAX+1
C     ADD ON NEW MASS WITH DENSITY=Z(111) IN TARGET
      DO 50 I=1,NIMAX
      K=(JJ-1)*IMAX+I+1
      DO 60 J=JJ,JMAX
      AMX(K)=Z(111)*TAU(I)*DY(J)
60    K=K+IMAX
50    CONTINUE
      JJ=(Z(147)/2.+2)
      JJ=JJ+1
      DO 61 I=11,IMAX
      K=I+1+(JJ-1)*IMAX
      DO 62 J=JJ,JMAX
      AMX(K)=Z(111)*TAU(I)*DY(J)
62    K=K+IMAX
61    CONTINUE
C     RESET ACTIVE GRID MARKERS.
      JJ=JJ-1
      Z(147)=JJ
      I1=NIMAX+2
      I2=NJMAX+2
      MS=I+DTNA
      NK=NC+1
C     EDIT THE NEW GRID.
      WRITE (6,8004)MS,NK,DX(1)
      WRITE (6,8007)IMAX,(X(I),I=0,IMAX)
      WRITE (6,8006)JMAX,(Y(J),J=0,JMAX)
      WRITE (6,8009)IMAX,(DX(I),I=1,IMAX)
      WRITE (6,8010)JMAX,(DY(J),J=1,JMAX)
      WRITE (6,8011)IMAX,(TAU(I),I=1,IMAX)
      KMAX=IMAX*JMAX+1
      IMAXA=IMAX+1
      JMAXA=JMAX+1
      KMAXA=KMAX+1
      RETURN
80040FORMAT(1H ////22H PROBLEM REZONED AT T=1PE12.6,6X,5HCYCLE14,6X,3HOREZONED
      1X=E12.6////)
8007 FORMAT(1H /10H X(I) I=0,12/(5F16.6))
8008 FORMAT(1H /10H Y(J) J=0,12/(5F16.6))
8009 FORMAT(1H /11H DX(I) I=1,12/(5F16.6))
8010 FORMAT(1H /11H DY(J) J=1,12/(5F16.6))
8011 FORMAT(1H /13H AREA(I) I=1,12/(F16.6,4F18.6))
      END

```

REZ02040
REZ02050
REZ02060
REZ02070

REZ02080
REZ02090
REZ02100
REZ02110
REZ02120
REZ02130
REZ02140
REZ02150
REZ02160
REZ02170
REZ02180

REZ02190
REZ02200
REZ02210
REZ02220
REZ02230
REZ02240

REZ02260

12.
HIBFTC EDIT LIST,DECK,REF
SUBROUTINE EDIT

EDIT0010
EDIT0490
EDIT1000

SENSE LITE (1) INDICATES LAST CYCLE OF THIS
RUN.
SENSE LITE (3) INDICATES FIRST CYCLE OF THIS
RUN.

104 CALL SLITET(3,K000FX)
GO TO(106,108),K000FX

EDIT1040
EDIT1050
EDIT1060
EDIT1070
EDIT1080

106 CALL SLITE (3)
GO TO 126

108 IF(CYCLE-CSTOP)110,122,122

110 IF(IREZ)9901,112,124

112 IF(AMOD(CYCLE,DUMPT7))114,124,114

EDIT1100

114 IF(AMOD(CYCLE,PRINTL))120,126,120

120 IF(AMOD(CYCLE,PRINTS))140,128,140

EDIT1150

NORMAL STOP ON THIS CYCLE.

122 CALL SLITE (1)

EDIT1160

DUMP ON TAPE 7.

124 GO TO 1

EDIT1170

126 CALL SLITE (4)

EDIT1180

128 GO TO 6000

EDIT1190

130 GO TO 1000

EDIT1200

132 CALL SLITET(4,K000FX)
GO TO(134,136),K000FX

EDIT1210

134 GO TO 5000

EDIT1220
EDIT1230

CHECK FOR ENERGY CHECK ERROR. WHERE

ECK= PERCENT ERROR/PER CYCLE.

$ECK = (ETH - E1) / ETH$ AT CYCLE $N - (ETH - E1) / ETH$

AT CYCLE $N - NPC$ ALL DIVIDED BY NPC, NOTE

$NPC = NO. OF CYCLES BETWEEN ENERGY CHECK$

136 IF(ABS(ECK)-DMIN)140,140,9905

EDIT1240

140 CALL SLITET(1,K000FX)
GO TO(142,144),K000FX

EDIT1250

EDIT1260

142 REWIND 7

CALL SLITE (1)

EDIT1280

144 GO TO 10000

EDIT1290

EDIT1300

EDIT1310

EDIT1320

EDIT1330

DUMP ON TAPE 7

1 IF(DUMPT7)30,3,3

3 BACKSPACE 7

REWIND 2

REWIND 3

MS=555.0

EDIT1360

WRITE (7)MS,CYCLE,N3

WRITE (7)(Z(L),L=1,MZ)

6 WRITE (7)(U(K),V(K),AMX(K),AIX(K),P(K),K=1,KMAX)

7 WRITE (7)(X(O),X(K),TAU(K),K=1,IMAX)

WRITE (7) Y(K), K=0, JMAX)	
C AGAIN, IF PROBLEM NO. IS NEGATIVE, WRITE	
C PARTICLE RECORDS ON TAPE 7.	
IF (PROB) 16, 16, 4	EDIT1420
16 DO 13 I=1, N3	EDIT1430
IF (N1) 150, 148, 150	
148 CONTINUE	
READ (2) (AM(N), XL(N), YL(N), I=1(N), I=2(N), N=2, N4)	
GO TO 152	
150 CONTINUE	
READ (3) (AM(N), XL(N), YL(N), I=1(N), I=2(N), N=2, N4)	
152 CONTINUE	
WRITE (7) (AM(N), XL(N), YL(N), I=1(N), I=2(N), N=2, N4)	
18 CONTINUE	EDIT1460
4 REWIND 2	
REWIND 3	
WS=666.0	EDIT1480
WRITE (7) WS, WS, WS	
WRITE (6, 8120) NC	EDIT1500
30 GO TO 126	EDIT1510
C	EDIT1520
C	EDIT1530
6000 NK=12	EDIT1540
C TABS ARE TANGENT ALPHAS.	
TAB(1)=0.02	EDIT1550
TAB(2)=0.04	EDIT1560
TAB(3)=0.06	EDIT1570
TAB(4)=0.08	EDIT1580
TAB(5)=0.10	EDIT1590
TAB(6)=0.15	EDIT1600
TAB(7)=0.20	EDIT1610
TAB(8)=0.25	EDIT1620
TAB(9)=0.30	EDIT1630
TAB(10)=0.4	EDIT1640
TAB(11)=0.5	EDIT1650
TAB(12)=1.0	EDIT1660
6010 DO 6012 I=1, 18	EDIT1670
6012 PR(I)=0.0	EDIT1680
NK1=NK+2	EDIT1690
DO 6014 I=1, NK1	EDIT1700
AMK(I)=0.0	EDIT1710
PK(I)=0.0	EDIT1720
6014 OK(I)=0.0	EDIT1730
DO 6028 K=2, KMAX	EDIT1740
6017 PR(1)=0.0	EDIT1760
PR(2)=0.0	EDIT1770
PR(4)=0.	
C CALCULATE KINETIC ENERGY IN CELL K.	
WS8=(U(K)**2+V(K)**2)*.5	
6019 IF (AMX(K)) 19917, 6028, 6020	EDIT1790

-34-

5020	I=MX1	EDIT1800
	IF(V(K))6026,6026,5022	EDIT1810
6022	WSA=ABS(U(K))/V(K)	EDIT1820
	DO 6024 I=1, MX	EDIT1830
C	SEARCH FOR TAN ANGLE THAT VELOCITY VECTORS	
C	MAKE.	
	IF(TAN(I)-WSA)6024,6026,6026	EDIT1840
6024	CONTINUE	EDIT1850
	I=MX+1	EDIT1860
6026	WS=AMX(K)	EDIT1870
C	SUM UP MASS BETWEEN ANGLES.	
6027	AMX(I)=AMX(I)+AMX(K)	EDIT1880
C	SUM UP RADIAL MOMENTA IN THE ANGLES.	
	PR(I)=PR(I)+U(K)*AMX(K)	EDIT1890
C	SUM UP AXIAL MOMENTA IN THE ANGLES.	
	QK(I)=QK(I)+V(K)*AMX(K)	EDIT1900
C	SUM UP TOTAL INTERNAL ENERGY	
	PR(5)=PR(5)+AIX(K)*AMX(K)	EDIT1910
C	SUM UP TOTAL KINETIC ENERGY	
	PR(6)=PR(6)+WSB*AMX(K)	EDIT1920
C	SUM UP TOTAL MASS	
	PR(8)=PR(8)+AMX(K)	EDIT1930
6028	CONTINUE	EDIT1940
	PR(3)=PR(1)+PR(2)	EDIT1950
	PR(7)=PR(5)+PR(6)	EDIT1960
	XNRG=PR(7)	EDIT1970
	PR(9)=PR(1)+PR(5)	EDIT1980
	PR(10)=PR(2)+PR(6)	EDIT1990
	PR(11)=PR(3)+PR(7)	EDIT2000
	PR(12)=PR(4)+PR(8)	EDIT2010
	WSA=(ETH-PR(11))/ETH	EDIT2020
	IF(CYCLE)9931,9931,9932	
9931	NPC=1	
9932	PR(18)=(WSA-DNA)/FLOAT(NPC)	
	ECK=PR(18)	EDIT2040
	DNA=WSA	EDIT2050
C	RESET CYCLE COUNTER BETWEEN ENERGY CHECK.	
	NPC=0	EDIT2060
	SUM=0.0	EDIT2070
	DO 800 I=1,13	EDIT2080
	SUM=SUM+QK(I)	EDIT2090
800	CONTINUE	EDIT2100
C	RADET= TOTAL POSITIVE AXIAL MOMENTUM IN GRID	
	RADET=SUM	EDIT2110
801	SUM=0.0	EDIT2120
	DO 810 K=2,KMAX	EDIT2130
	IF(AMX(K))810,810,802	EDIT2140
802	IF(U(K))810,810,803	EDIT2150
803	SUM=SUM+AMX(K)*U(K)	EDIT2160
810	CONTINUE	EDIT2170

[illegible]

136.

1160	PR(I)=OR(PR(I), ABS(922746880))	EDIT2650
	GO TO 1180	EDIT2660
C	REPLACE 6000000000 BY 805306368	EDIT2670
1166	PR(I)=OR(PR(I), ABS(805306368))	EDIT2680
1180	CONTINUE	EDIT2690
1200	IF(MOD(J,5))1210,1204,1210	EDIT2700
1204	IF(DY(J)-DY(J-1))1206,1208,1206	EDIT2710
1206	WRITE (6,8211)DY(J),J,(PR(I),I=1,I1)	EDIT2720
	GO TO 1224	EDIT2730
1208	WRITE (6,8201)J,(PR(I),I=1,I1)	EDIT2740
	GO TO 1224	EDIT2750
1210	IF(DY(J)-DY(J-1))1212,1214,1212	EDIT2760
1212	WRITE (6,8222)DY(J),(PR(I),I=1,I1)	EDIT2770
	GO TO 1224	EDIT2780
1214	WRITE (6,8202)(PR(I),I=1,I1)	EDIT2790
1224	J=J-1	EDIT2800
1226	IF(J)1230,1230,1100	EDIT2810
C	REPLACE 604000000000 BY-17716740096	EDIT2820
1230	PR(I)=(-ABS(-17716740096))	EDIT2830
	WRITE (6,8201)J,(PR(I),I=1,I1)	EDIT2840
	WRITE (6,8302)(I,I=0,IMAX,5)	EDIT2850
1240	GO TO 132	EDIT2860
C****	END OF PLOT SUBROUTINE *****	EDIT2870
C		EDIT2880
C		EDIT2890
C****	SUBROUTINE L P *****	EDIT2900
5000	WRITE (6,8116)PROB,NC,T,DTNA,TRAD,DTRAD,NR,N1,N2,N3,N4	EDIT2910
5004	DO 5050 I=1,I1	EDIT2920
	CALL SLITE (4)	EDIT2930
	J=I2+1	EDIT2940
	K=I2*IMAX+1+I	EDIT2950
	DO 5046 L=1,I2	EDIT2960
	J=J-1	EDIT2970
	K=K-IMAX	EDIT2980
5012	IF(AMX(K))9917,5046,5014	EDIT2990
5014	CALL SLITET(4,K000FX)	EDIT3000
	GO TO(5016,5019),K000FX	EDIT3010
5016	WRITE (6,8135)I,X(I),DX(I)	EDIT3020
C	WS= DENSITY OF CELL(K) IN GRAMS/CM. CUBED.	
5019	WS=AMX(K)/((TAU(I)*DY(J))	EDIT3030
C	WSA= COMPRESSION = RHO/RHO NOT.	
	WSA=WS/Z(111)	EDIT3040
C	WSC= PRESSURE CONVERTED TO MEGABARS.	
	WSC=P(K)*1.E+4	EDIT3050
C	FIRST COLUMN= (J) THE ROW NO.	
C	SECOND COLUMN= RADIAL VELOCITY CM./SHAKE	
C	THIRD COLUMN= AXIAL VELOCITY CM./SHAKE	
C	FOURTH COLUMN= F/A = PRESSURE IN MEGABARS	
C	FIFTH COLUMN = AMX = MASS IN GRAMS.	
C	SIXTH COLUMN = RHO = DENSITY IN GRAMS/CC.	

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C SEVENTH COLUMN = AIX = SPECIFIC INTERNAL ENERGY JERKS/GM.
C EIGHT COLUMN = COMPRESSION = RHO/RHO NOT
C NINTH COLUMN = Z VALUE (CM.) OF TOP OF CELL)
50180WRITE (6,8108)J,U(K),V(K),WSC,AMX(K), WS,AIX(K),EDIT3060
      1WSA,Y(J) EDIT3070
5046 CONTINUE EDIT3080
5050 CONTINUE EDIT3090
      GO TO 136 EDIT3100
C**** END OF L P SUBROUTINE ***** EDIT3110
C EDIT3120
C EDIT3130
C ERROR EDIT3140
9901 NK=110 EDIT3150
      GO TO 9999 EDIT3160
C ENERGY CHECK EDIT3170
9905 NK=136 EDIT3180
      GO TO 9999 EDIT3190
C NEGATIVE MASS EDIT3200
9917 NK=6015 EDIT3210
      GO TO 9999 EDIT3220
9920 NK=904 EDIT3230
      GO TO 9999 EDIT3240
9921 NK=912 EDIT3250
      GO TO 9999 EDIT3260
9922 NK=918 EDIT3270
      GO TO 9999 EDIT3280
9923 NK=922 EDIT3290
      GO TO 9999 EDIT3300
9924 NK=926 EDIT3310
9999 NR=6 EDIT3320
      CALL DUMP EDIT3330
10000 RETURN EDIT3340
C EDIT3350
C FORMATS EDIT3360
8108 FORMAT(I3,1X,1P2E14.6,3E15.6,E14.6,E15.6,E14.6) EDIT3370
81160FORMAT(8H1PROBLEM6X,5HCYCLE9X,4HTIME13X,2HDT13X,4HTRAD11X,5HDTRAD1EDIT3380
      12X,2HNR6X,2HN14X,2HN24X,2HN34X,2HN4/(F7.1,I11,2X,1P4E16.7,I10,2X,4EDIT3390
      2I6)) EDIT3400
81170FORMAT(1H0//17X2HAI16X,2HAK14X,5HAI+AK15X,2HAM/4H DOT3X,1P4E18.7/3EDIT3410
      1H X4X,4E18.7) EDIT3420
81180FORMAT(12X,13H-----5X,13H-----5X,13H-----5EDIT3430
      1X,13H-----/7H TOTALS1P4E18.7) EDIT3440
81190FORMAT(2H0 //16X,5HRADEB13X,5HRADEB13X,5HRADET12X,7HMAX VEL13X,3HTEDIT3450
      1HE12X,9HREL ERROR/7X,1P6E18.7////) EDIT3460
8120 FORMAT(1H0//21H TAPE 7 DUMP ON CYCLE15////) EDIT3470
81240FORMAT(3H K12X,5HAM(K)11X,9HSUM AM(K)11X,4HP(K)13X,4HQ(K)/(I3,4X,EDIT3480
      11P4E18.7)) EDIT3490
8131 FORMAT(1H ///11H DY(J) J=1,I2//(10F12.3)) EDIT3500
8133 FORMAT(1H ///11H Y(J) J=0,I2//(10F12.3)) EDIT3510
81350FORMAT(1H ///4H I =I3,6X,6HX(I) =F12.3,6X,7HDX(I) =F12.3//3H J8X,EDIT3520

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138.

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      11HX13X,1HY13X,3HF/A12X,3HAMX12X,3HRHO11X,3HAI1X12X,4HCOMP11X,2H Y/) EDIT3530
8201 FORMAT(110,2H I54A2) EDIT3540
8202 FORMAT(10X,2H I54A2) EDIT3550
8211 FORMAT(F7.1,I3,2H I54A2) EDIT3560
8222 FORMAT(F7.1,3X,2H I54A2) EDIT3570
8302 FORMAT(112,10I10) EDIT3580
83070FORMAT(5H X1 =1PE12.6,3X,4HX2 =E12.6,3X,6HXMAX =E12.6,6X,4HY1 =E12 EDIT3590
      1.6,3X,4HY2 =E12.6,3X,6HYMAX =E12.6) EDIT3600
8308 FORMAT(1H /) EDIT3610
9040 FORMAT(1H / 616) EDIT3620
      END EDIT3630
```

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